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For

: 2,3,6-TRISUBSTITUTED-4-PYRIMIDONE DERIVATIVES

# **CLAIM OF PRIORITY**

Commissioner for Patents U.S. Patent and Trademark Office Customer Service Window, Mail Stop PCT Randolph Building 401 Dulany Street Alexandria, VA 22314

Sir:

Applicant hereby claims the right of priority granted pursuant to 35 U.S.C. 119 and 365 based upon Japanese Application Nos. 2003-126021, filed March 26, 2003 and 2003-126022, filed March 26, 2003. The International Bureau already should have sent certified copies of the Japanese applications to the United Stated designated office. If the certified copies have not arrived, please contact the undersigned.

> Respectfully submitted, Kazutoshi WATANABE et al.

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# 日本国特許庁 JAPAN PATENT OFFICE

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別紙添付の書類に記載されている事項は下記の出願書類に記載されている事項と同一であることを証明する。

This is to certify that the annexed is a true copy of the following application as filed with this Office.

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Applicant(s):

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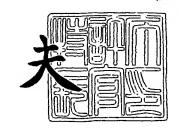
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【その他】 サノフィーサンテラボはフランス国の法律に基づく法人

【提出物件の目録】

【物件名】 外国語明細書 1

【物件名】 外国語要約書 1

1

# 【書類名】 外国語明細書

# 1. TITLE OF INVENTION

## 3-SUBSTITUTED-4-PYRIMIDONE DERIVATIVES

#### 2. CLAIMS

1. A pyrimidone derivative represented by formula (I) or a salt thereof, or a solvate thereof or a hydrate thereof:

$$(X)_{m} \xrightarrow{N}_{R} 0$$

$$(Y)_{0}$$

$$(Y)_{0}$$

$$(Y)_{0}$$

$$(Y)_{0}$$

$$(Y)_{0}$$

$$(Y)_{0}$$

$$(Y)_{0}$$

wherein R represents a C1-C12 alkyl group which may be substituted; the ring of:

represents piperazine ring or piperidine ring;

each X independently represents a C<sub>1</sub>-C<sub>8</sub> alkyl group which may be substituted, an optionally partially hydrogenated C<sub>6</sub>-C<sub>10</sub> aryl ring which may be substituted, an indan ring which may be substituted, an optionally substituted heterocyclic ring having 1 to 4 hetero atoms selected from the group consisting of oxygen atom, sulfur atom, and nitrogen atom, and having 5 to 10 ring-constituting atoms in total, a phenylamino group which may be substituted, or X¹-CO- wherein X¹ represents a C<sub>6</sub>-C<sub>10</sub> aryl ring which may be substituted or a C<sub>1</sub>-C<sub>8</sub> alkyl group which may be substituted;

m represents an integer of 1 to 3;

each Y independently represents a halogen atom, a hydroxy group, a cyano group, a  $C_1$ - $C_6$  alkyl group which may be substituted, Y¹-CO- wherein Y¹ represents a  $C_6$ - $C_{10}$  aryl ring which may be substituted or a  $C_1$ - $C_8$  alkyl group which may be substituted or  $Y^2$ -O-CO- wherein  $Y^2$  represents a  $C_1$ - $C_8$  alkyl group which may be substituted;

n represents an integer of 0 to 8;

when X and Y or two Y groups are attached on the same carbon atom, they may combine to each other to form a  $C_2$ - $C_6$  alkylene group; and when m is 1, n is 0, and X is  $X^1$ -CO-,

- (1) X does not bind to 3-position of unsubstituted 1-piperazinyl group or does not bind to 3-position of a 4-alkyl-1-piperazinyl group; or
- (2) X does not bind to 3-position or 4-position of non-substituted 1-piperidinyl group.
- 2. The pyrimidone derivative or the salt thereof, or the solvate thereof or the hydrate thereof according to claim 1 having the following formula(II)

$$(X)_{p} \qquad (Y)_{r} \qquad (II)$$

$$(X)_{q} \qquad (Y)_{r} \qquad (X)_{q} \qquad (Y)_{r} \qquad (X)_{q} \qquad (Y)_{r} \qquad (X)_{q} \qquad (X)_{q}$$

wherein R, X and Y are the same as those defined in claim 1; p is 0 or 1; q is 0 or 1; r is an integer of 0 to 6; p+q is 1 or 2; and Z represents N or CZ<sup>1</sup> wherein Z<sup>1</sup> represents hydrogen atom or Y.

- 3. The pyrimidone derivative or the salt thereof, or the solvate thereof or the hydrate thereof according to claim 2, wherein R is a C<sub>1</sub>-C<sub>3</sub> alkyl group which may be substituted by a C<sub>3</sub>-C<sub>3</sub> cycloalkyl group.
- 4. The pyrimidone derivative or the salt thereof, or the solvate thereof or the hydrate thereof according to claim 3, wherein R is methyl group or ethyl group; Y is in 3-, 4- or 5-position of the piperazine ring or the piperidine ring; p+q is 1; and r is an integer of 0 to 3.
- 5. The pyrimidone derivative or the salt thereof, or the solvate thereof or the hydrate thereof according to claim 4, wherein X is a C<sub>1</sub>-C<sub>8</sub> alkyl group which may be substituted or a C<sub>6</sub>-C<sub>10</sub> aryl ring which may be substituted; Y is a C<sub>1</sub>-C<sub>6</sub> alkyl group which may be substituted; p is 1; q is 0; r is an integer of 0 to 3; and Z is N or

CH.

one;

- 6. The pyrimidone derivative or the salt thereof, or the solvate thereof or the hydrate thereof according to claim 5, wherein X is a benzene ring which may be substituted, a benzyl group which may be substituted; Y is a methyl group which may be substituted; Z is N and r is 0 or 1.
- 7. The pyrimidone derivative or the salt thereof, or the solvate thereof or the hydrate thereof according to claim 4, wherein X is a benzene ring which may be substituted, a benzyl group which may be substituted, a benzoyl group which may be substituted; Y is a methyl group which may be substituted; Y is a methyl group which may be substituted; Z is N and p is 0.
- 8. The pyrimidone derivative or the salt thereof, or the solvate thereof or the hydrate thereof according to claim 4, wherein X is a  $C_1$ - $C_8$  alkyl group substituted by a benzene ring which may be substituted or a benzene ring which may be substituted; Y is a hydroxy group, a cyano group, or Y¹-CO- wherein Y¹ is a  $C_1$ - $C_8$  alkyl group; Z is CH or C-Y and r is 0 or 1.
- 9. The pyrimidone derivative or the salt thereof, or the solvate thereof or the hydrate thereof according to claim 8, wherein X is a benzyl group which may be substituted or a benzene ring which may be substituted; Y is a hydroxy group, a cyano group, or an acetyl group; Z is CH or C-Y and r is 0 or 1.
- 10. A pyrimidone derivative which is selected from the group consisting of: 2-(3-Phenylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; 2-(3-(4-Fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; 2-(3-(3-Fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; 2-(3-(2-Fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; (S)-2-(3-(4-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- (R)-2-(3-(4-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(3-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- $2\hbox{-}(3\hbox{-}(4\hbox{-Bromophenyl}) piperazin-1\hbox{-}yl)-3\hbox{-methyl-}6\hbox{-}(4\hbox{-pyridyl})-3\textit{H-pyrimidin-4-one};$

- 2-(3-(3-Bromophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Bromophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(4-Methylphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(3-Methylphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Methylphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(4-Cyanophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(3-Cyanophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Cyanophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(4-Methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(3-Methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Ethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(5-Fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one; 2-(3-(4-Fluoro-3-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-
- pyrimidin-4-one;
- 2-(3-(4-Fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;
- (S)-2-(3-(4-Fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;
- (R)-2-(3-(4-Fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;
- 2-(3-(4-Chloro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;
- 2-(3-(4-Fluoro-2-methylphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;
- 2-(3-(2-Fluoro-6-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;
- 2-(3-(5-Bromo-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;
- 2-(3-(2-Bromo-4-fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;
- 2-(3-(2-Chloro-6-fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one;

- 2-(3-(2,4-Difluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(2,6-Difluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(2,6-Dichlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2,4-Dimethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(3,4-Dimethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(2,5-Dimethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(2,6-Dimethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one:
- 2-(3-(2,4-Difluoro-6-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;(1034)
- 2-(3-(5-Cyano-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-Cyano-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(1-Naphthyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2-Naphthyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2,3-Dihydrobenzofuran-7-yl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(Benzofuran-2-yl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- (S)-2-(3-(Benzofuran-2-yl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-(Pyrrolidin-1-yl-methyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-(Pyrrolidin-1-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(2-methoxy-4-(pyrrolidin-1-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(2-methoxy-5-(pyrrolidin-1-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-

- 3 H-pyrimidin-4-one;
- 2-(3-(4-(Phenyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-(4-Fluorophenyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(4-(4-Methoxyphenyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-(2-Methoxyphenyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-(Morpholin-4-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-(4-Methylpiperazin-1-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(4-Phenylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(4-Benzylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(4-Benzoylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(4-(1,2-Benzisothiazol-3-yl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(4-Methyl-3-phenylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-Fluoro-2-methoxyphenyl)-4-methylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3-H-pyrimidin-4-one;
- (S)-2-(3-(4-Fluoro-2-methoxyphenyl)-4-methylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- (R)-2-(3-(4-Fluoro-2-methoxyphenyl)-4-methylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(4-Acetyl-3-(4-fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-
- 3 H-pyrimidin-4-one;
- $2\hbox{-}(4\hbox{-}Benzyl\hbox{-}3\hbox{-}(4\hbox{-}fluoro\hbox{-}2\hbox{-}methoxyphenyl) piperazin\hbox{-}1\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}10\hbox{-}yl)\hbox{-}3\hbox{-}methyl\hbox{-}6\hbox{-}(4\hbox{-}pyridyl)\hbox{-}3\hbox{-}methyl]$
- 3 H-pyrimidin-4-one;
- 2-(4-Benzyl-3-(ethoxycarbonyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(4-methyl-3-(1-naphthyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(5,5-Dimethyl-3-(2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;

- 2-(3-(4-Fluorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(3-Fluorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(3-Fluorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(4-Chlorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(4-Chlorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(4-Bromophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(4-Methoxyphenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(4-Methoxyphenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(4-((-1-y)-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
  2-(3-(4-((-1-y)-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- (S)-2-(3-(4-(Pyrrolidin-1-yl-methyl)phenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- (R)-2-(3-(4-(Pyrrolidin-1-yl-methyl)phenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; and
- 2-(3-Hydroxy-3-phenylpiperidin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one, or a salt thereof, or a solvate thereof or a hydrate thereof.
- 11. A medicament comprising as an active ingredient a substance selected from the group consisting of the pyrimidone derivative represented by formula (I) and a salt thereof, and a solvate thereof and a hydrate thereof according to claim 1.
- 12. A tau protein kinase 1 inhibitor selected from the group consisting of the pyrimidone derivative represented by formula (I) and a salt thereof, and a solvate thereof and a hydrate thereof according to claim 1.
- 13. The medicament according to claim 11 which is used for preventive and/or therapeutic treatment of a disease caused by tau protein kinase 1 hyperactivity.
- 14. The medicament according to claim 11 which is used for preventive and/or therapeutic treatment of a neurodegenerative disease.
- 15. The medicament according to claim 14, wherein the neurodegenerative disease is selected from the group consisting of Alzheimer disease, ischemic cerebrovascular accidents, Down syndrome, cerebral bleeding due to cerebral amyloid angiopathy, progressive supranuclear palsy, subacute sclerosing panencephalitic parkinsonism, postencephalitic parkinsonism, pugilistic

encephalitis, Guam parkinsonism-dementia complex, Lewy body disease, Pick's disease, corticobasal degeneration, frontotemporal dementia, vascular dementia, traumatic injuries, brain and spinal cord trauma, peripheral neuropathies, retinopathies, and glaucoma.

16. The medicament according to claim 11, wherein the disease is selected from the group consisting of non-insulin dependent diabetes, obesity, manic depressive illness, schizophrenia, alopecia, breast cancer, non-small cell lung carcinoma, thyroid cancer, T or B-cell leukemia, and a virus-induced tumor.

# 3. DETAILED DESCRIPTION OF INVENTION Technical Field

The present invention relates to compounds that are useful as an active ingredient of a medicament for preventive and/or therapeutic treatment of diseases mainly caused by abnormal activity of tau protein kinase 1, such as neurodegenerative diseases (e.g. Alzheimer disease).

# Background Art

Alzheimer disease is progressive senile dementia, in which marked cerebral cortical atrophy is observed due to degeneration of nerve cells and decrease of nerve cell number. Pathologically, numerous senile plaques and neurofibrillary tangles are observed in brain. The number of patients has been increased with the increment of aged population, and the disease arises a serious social problem. Although various theories have been proposed, a cause of the disease has not yet been elucidated. Early resolution of the cause has been desired.

It has been known that the degree of appearance of two characteristic pathological changes of Alzheimer disease well correlates to the degree of intellectual dysfunction. Therefore, researches have been conducted from early 1980's to reveal the cause of the disease through molecular level investigations of components of the two pathological changes. Senile plaques accumulate extracellularly, and  $\beta$  amyloid protein has been elucidated as their main component (abbreviated as "A  $\beta$ " hereinafter in the specification: Biochem. Biophys. Res. Commun., 120, 855 (1984); EMBO J., 4, 2757 (1985); Proc. Natl. Acad. Sci. USA,

82, 4245 (1985)). In the other pathological change, i.e., the neurofibrillary tangles, a double-helical filamentous substance called paired helical filament (abbreviated as "PHF" hereinafter in the specification) accumulate intracellularly, and tau protein, which is a kind of microtubule-associated protein specific for brain, has been revealed as its main component (Proc. Natl. Acad. Sci. USA, 85, 4506 (1988); Neuron, 1, 827 (1988)).

Furthermore, on the basis of genetic investigations, presenilins 1 and 2 were found as causative genes of familial Alzheimer disease (Nature, 375, 754 (1995); Science, 269, 973 (1995); Nature. 376, 775 (1995)), and it has been revealed that presence of mutants of presenilins 1 and 2 promotes the secretion of A  $\beta$  (Neuron, 17, 1005 (1996); Proc. Natl. Acad. Sci. USA, 94, 2025 (1997)). From these results, it is considered that, in Alzheimer disease, A  $\beta$  abnormally accumulates and agglomerates due to a certain reason, which engages with the formation of PHF to cause death of nerve cells. It is also expected that extracellular outflow of glutamic acid and activation of glutamate receptor responding to the outflow may possibly be important factors in an early process of the nerve cell death caused by ischemic cerebrovascular accidents (Sai-shin Igaku [Latest Medicine], 49, 1506 (1994)).

It has been reported that kainic acid treatment that stimulates the AMPA receptor, one of glutamate receptor, increases mRNA of the amyloid precursor protein (abbreviated as "APP" hereinafter in the specification) as a precursor of A  $\beta$  (Society for Neuroscience Abstracts, 17, 1445 (1991)), and also promotes metabolism of APP (The Journal of Neuroscience, 10, 2400 (1990)). Therefore, it has been strongly suggested that the accumulation of A $\beta$  is involved in cellular death due to ischemic cerebrovascular disorders. Other diseases in which abnormal accumulation and agglomeration of A $\beta$  are observed include, for example, Down syndrome, cerebral bleeding due to solitary cerebral amyloid angiopathy, Lewy body disease (Shin-kei Shinpo [Nerve Advance], 34, 343 (1990); Tanpaku-shitu Kaku-san Koso [Protein, Nucleic Acid, Enzyme], 41, 1476 (1996)) and the like. Furthermore, as diseases showing neurofibrillary tangles due to the PHF accumulation, examples include progressive supranuclear palsy, subacute sclerosing panencephalitic parkinsonism, postencephalitic parkinsonism, pugilistic encephalitis, Guam

parkinsonism-dementia complex, Lewy body disease and the like (Tanpakushitu Kakusan Koso [Protein, Nucleic Acid, Enzyme], 36, 2 (1991); Igaku no Ayumi [Progress of Medicine], 158, 511 (1991); Tanpakushitu Kakusan Koso [Protein, Nucleic Acid, Enzyme], 41, 1476 (1996)).

The tau protein is generally composed of a group of related proteins that forms several bands at molecular weights of 48-65 kDa in SDS-polyacrylamide gel electrophoresis, and it promotes the formation of microtubules. It has been verified that tau protein incorporated in the PHF in the brain suffering from Alzheimer disease is abnormally phosphorylated compared with usual tau protein (J. Biochem., 99, 1807 (1986); Proc. Natl. Acad. Sci. USA, 83, 4913 (1986)). An enzyme catalyzing the abnormal phosphorylation has been isolated. The protein was named as tau protein kinase 1 (abbreviated as "TPK1" hereinafter in the specification), and its physicochemical properties have been elucidated (Seikagaku [Biochemistry], 64, 308 (1992); J. Biol. Chem., 267, 10897 (1992)). Moreover, cDNA of rat TPK1 was cloned from a rat cerebral cortex cDNA library based on a partial amino acid sequence of TPK1, and its nucleotide sequence was determined and an amino acid sequence was deduced (Japanese Patent Un-examined Publication [Kokai] No. 6-239893/1994). As a result, it has been revealed that the primary structure of the rat TPK1 corresponds to that of the enzyme known as rat GSK-3  $\beta$  (glycogen synthase kinase  $3\beta$ , FEBS Lett., 325, 167 (1993)).

It has been reported that A  $\beta$ , the main component of senile plaques, is neurotoxic (Science, 250, 279 (1990)). However, various theories have been proposed as for the reason why A  $\beta$  causes the cell death, and any authentic theory has not yet been established. Takashima et al. observed that the cell death was caused by A  $\beta$  treatment of fetal rat hippocampus primary culture system, and then found that the TPK1 activity was increased by A  $\beta$  treatment and the cell death by A  $\beta$  was inhibited by antisense of TPK1 (Proc. Natl. Acad. Sci. USA, 90, 7789 (1993); Japanese Patent Un-examined Publication [Kokai] No. 6-329551/1994).

In view of the foregoing, compounds which inhibit the TPK1 activity may possibly suppress the neurotoxicity of A  $\beta$  and the formation of PHF and inhibit the nerve cell death in the Alzheimer disease, thereby cease or defer the progress of the disease. The compounds may also be possibly used as a medicament for therapeutic

treatment of ischemic cerebrovascular disorder, Down syndrome, cerebral amyloid angiopathy, cerebral bleeding due to Lewy body disease and the like by suppressing the cytotoxicity of A  $\beta$ . Furthermore, the compounds may possibly be used as a medicament for therapeutic treatment of neurodegenerative diseases such as progressive supranuclear palsy, subacute sclerosing panencephalitic parkinsonism, postencephalitic parkinsonism, pugilistic encephalitis, Guam parkinsonism-dementia complex, Lewy body disease, Pick's disease, corticobasal degeneration, frontotemporal dementia, vascular dementia, acute stroke and traumatic injuries, brain and spinal cord trauma, peripheral neuropathies, retinopathies and glaucoma; non-insulin dependent diabetes (such as diabetes type II), and obesity, manic depressive illness, schizophrenia, alopecia, cancers such as breast cancer, non-small cell lung carcinoma, thyroid cancer, T or B-cell leukemia and several virus-induced tumors.

As structurally similar compounds to the compounds of the present invention represented by formula (I) described later, compounds represented by the following formula (A) are known:

wherein R represents 2,6-dichlorobenzyl group, 2-(2-chlorophenyl)ethylamino group, 3-phenylpropylamino group, or 1-methyl-3-phenylpropylamino group (W098/24782). The compounds represented by formula (A) are characterized to have 4-fluorophenyl group at the 5-position of the pyrimidine ring and a hydroxy group at the 4-position, and not falling within the scope of the present invention. Moreover, main pharmacological activity of the compounds represented by formula (A) is anti-inflammatory effect, whereas the compounds of the present invention represented by formula (I) are useful as a TPK1 inhibitor or a medicament for

therapeutic treatment of neurodegenerative diseases, and therefore, their pharmacological activities are totally different to each other.

Patent Document 1: WO 00/18758

Patent Document 2: WO 01/70728

Patent Document 3: WO 01/70729

## Object to be Achieved by the Invention

An object of the present invention is to provide compounds useful as an active ingredient of a medicament for preventive and/or therapeutic treatment of diseases such as Alzheimer disease. More specifically, the object is to provide novel compounds useful as an active ingredient of a medicament that enables radical prevention and/or treatment of the neurodegenerative diseases such as Alzheimer disease by inhibiting the TPK1 activity to suppress the neurotoxicity of A  $\beta$  and the formation of the PHF and by inhibiting the death of nerve cells.

#### Means to Achieve the Object

In order to achieve the foregoing object, the inventors of the present invention conducted screenings of various compounds having inhibitory activity against the phosphorylation of TPK1. As a result, they found that compounds represented by the following formula (I) had the desired activity and were useful as an active ingredient of a medicament for preventive and/or therapeutic treatment of the aforementioned diseases. The present invention was achieved on the basis of these findings.

The present invention thus provides 3-substituted-4-pyrimidone derivatives represented by formula (I) or salts thereof, or solvates thereof or hydrates thereof:

$$(X)_{m} \longrightarrow N \longrightarrow N$$

$$(Y)_{n} \longrightarrow N \longrightarrow N$$

$$R$$

$$(I)$$

wherein R represents a C<sub>1</sub>-C<sub>12</sub> alkyl group which may be substituted; the ring of:

 $\binom{N}{2}$ 

represents piperazine ring or piperidine ring;

each X independently represents a C<sub>1</sub>-C<sub>8</sub> alkyl group which may be substituted, an optionally partially hydrogenated C<sub>6</sub>-C<sub>10</sub> aryl ring which may be substituted, an indan ring which may be substituted, an optionally substituted heterocyclic ring having 1 to 4 hetero atoms selected from the group consisting of oxygen atom, sulfur atom, and nitrogen atom, and having 5 to 10 ring-constituting atoms in total, a phenylamino group which may be substituted or X<sup>1</sup>-CO- wherein X<sup>1</sup> represents a C<sub>6</sub>-C<sub>10</sub> aryl ring which may be substituted or a C<sub>1</sub>-C<sub>8</sub> alkyl group which may be substituted;

m represents an integer of 1 to 3;

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each Y independently represents a halogen atom, a hydroxy group, a cyano group, a  $C_1$ - $C_6$  alkyl group which may be substituted, Y¹-CO- wherein Y¹ represents a  $C_6$ - $C_{10}$  aryl ring which may be substituted or a  $C_1$ - $C_8$  alkyl group which may be substituted or Y²-O-CO- wherein Y² represents a  $C_1$ - $C_8$  alkyl group which may be substituted; n represents an integer of 0 to 8;

when X and Y or two Y groups are attached on the same carbon atom, they may combine to each other to form a  $C_2$ - $C_6$  alkylene group; and when m is 1, n is 0, and X is  $X^1$ -CO-,

- (1) X does not bind to 3-position of unsubstituted 1-piperazinyl group or does not bind to 3-position of a 4-alkyl-1-piperazinyl group; or
- (2) X does not bind to 3-position or 4-position of non-substituted1-piperidinyl group.

According to another aspect of the present invention, there is provided a medicament comprising as an active ingredient a substance selected from the group consisting of the 3-substituted-4-pyrimidone derivatives represented by formula (I) and the physiologically acceptable salts thereof, and the solvates thereof and the hydrates thereof. As preferred embodiments of the medicament, there are provided the aforementioned medicament which is used for preventive and/or therapeutic treatment of diseases caused by tau protein kinase 1 hyperactivity, and the aforementioned medicament which is used for preventive and/or therapeutic treatment of neurodegenerative diseases.

As further preferred embodiments of the present invention, there are provided the aforementioned medicament wherein the diseases are selected from the group consisting of Alzheimer disease, ischemic cerebrovascular accidents, Down syndrome, cerebral bleeding due to cerebral amyloid angiopathy, progressive supranuclear palsy, subacute sclerosing panencephalitic parkinsonism, postencephalitic parkinsonism, pugilistic encephalitis, Guam parkinsonism-dementia complex, Lewy body disease, Pick's disease, corticobasal degeneration and frontotemporal dementia, vascular dementia, acute stroke and traumatic injuries, brain and spinal cord trauma, peripheral neuropathies, retinopathies and glaucoma, non-insulin dependent diabetes (such as diabetes type II), and obesity, manic depressive illness, schizophrenia, alopecia, cancers such as breast cancer, non-small cell lung carcinoma, thyroid cancer, T or B-cell leukemia and several virus-induced tumors; and the aforementioned medicament in the form of pharmaceutical composition containing the above substance as an active ingredient together with one or more pharmaceutical additives.

The present invention further provides an inhibitor of tau protein kinase 1 comprising as an active ingredient a substance selected from the group consisting of the 3-substituted-4-pyrimidone derivatives of formula (I) and the salts thereof, and the solvates thereof and the hydrates thereof.

According to further aspects of the present invention, there are provided a method for preventive and/or therapeutic treatment of diseases caused by tau protein kinase 1 hyperactivity, which comprises the step of administering to a

patient a preventively and/or therapeutically effective amount of a substance selected from the group consisting of the 3-substituted-4-pyrimidone derivatives of formula (I) and the physiologically acceptable salts thereof, and the solvates thereof and the hydrates thereof; and a use of a substance selected from the group consisting of the 3-substituted-4-pyrimidone derivatives of formula (I) and the physiologically acceptable salts thereof, and the solvates thereof and the hydrates thereof for the manufacture of the aforementioned medicament.

#### Mode for Carrying Out the Invention

The alkyl group used herein may be either linear or branched.

The C1-C12 alkyl group represented by R may be, for example, methyl group, ethyl group, n-propyl group, isopropyl group, n-butyl group, isobutyl group, sec-butyl group, tert-butyl group, n-pentyl group, isopentyl group, neopentyl group, 1,1-dimethylpropyl group, n-hexyl group, isohexyl group, or a linear or branched heptyl group, octyl group, nonyl group, decyl group, undecyl group or dodecyl group. Particularly preferred R is methyl group.

In the specification, when a functional group is defined as "which may be substituted" or "optionally substituted", the number of substituents as well as their types and substituting positions are not particularly limited, and when two or more substituents are present, they may be the same or different.

When the C<sub>1</sub>-C<sub>12</sub> alkyl group represented by R has one or more substituents, the alkyl group may have one or more substituents selected from, for example, the groups consisting of a C<sub>3</sub>-C<sub>8</sub> cycloalkyl group such as cyclopropyl group, cyclobutyl group, cyclopentyl group, cyclohexyl group, cycloheptyl group, cyclooctyl group; a C<sub>1</sub>-C<sub>5</sub> alkoxy group such as methoxy group, ethoxy group, propoxy group, isopropoxy group, butoxy group, isobutoxy group, tert-butoxy group; C<sub>1</sub>-C<sub>3</sub> alkylamino group or C<sub>2</sub>-C<sub>6</sub> dialkylamino group; a C<sub>6</sub>-C<sub>10</sub> aryl group such as phenyl group, 1-naphthyl group, and 2-naphthyl group.

The C1-C8 alkyl group represented by X may be, for example, methyl group, ethyl group, n-propyl group, isopropyl group, n-butyl group, isobutyl group, sec-butyl group, tert-butyl group, n-pentyl group, isopentyl group, neopentyl group, 1,1-dimethylpropyl group, n-hexyl group, isohexyl group, or a linear or branched

heptyl group or octyl group.

The optionally partially hydrogenated  $C_6$ - $C_{10}$  aryl ring represented by X may be, for example a benzene ring, a naphthalene ring, an indan ring or a 1,2,3,4-tetrahydronaphthalene ring.

The heterocyclic ring having 1 to 4 hetero atoms selected from the group consisting of oxygen atom, sulfur atom, and nitrogen atom, and having 5 to 10 ring-constituting atoms in total represented by X may be, for example, furan ring, dihydrofuran ring, tetrahydrofuran ring, pyran ring, dihydropyran ring, tetrahydropyran ring, benzofuran ring, dihydrobenzofuran, isobenzofuran ring, benzodioxol ring, chromene ring, chroman ring, isochroman ring, thiophene ring, benzothiophene ring, pyrrole ring, pyrroline ring, pyrrolidine ring, imidazole ring, imidazoline ring, imidazolidine ring, pyrazole ring, pyrazoline ring, pyrazolidine ring, triazole ring, tetrazole ring, pyridine ring, pyridine oxide ring, piperidine ring, pyrazine ring, piperazine ring, pyrimidine ring, pyridazine ring, indole ring, indoline ring, isoindole ring, isoindoline ring, indazole ring, benzimidazole ring, benzotriazole ring, tetrahydroisoquinoline ring, benzothiazolinone ring, benzoxazolinone ring, purine ring, quinolizine ring, quinoline ring, phthalazine ring, naphthyridine ring, quinoxaline ring, quinazoline ring, cinnoline ring, pteridine ring, oxazole ring, oxazolidine ring, isoxazole ring, isoxazolidine ring, oxadiazole ring, thiazole ring, benzothiazole ring, thiazylidine ring, isothiazole ring, isothiazolidine ring, benzodioxole ring, dioxane ring, benzodioxane ring, dithian ring, morpholine ring, thiomorpholine ring, or phthalimide ring.

The  $C_6$ - $C_{10}$  aryl ring represented by  $X^1$  or  $Y^1$  may be, for example, a benzene ring or a naphthalene ring.

When the ring represented by X or X¹ has one or more substituents, the ring may have one or more substituents selected form the group consisting of a C¹-C⁵ alkyl group such as methyl group, ethyl group, propyl group, isopropyl group, butyl group, isobutyl group, sec-butyl group, tert-butyl group, pentyl group, isopentyl group, neopentyl group, 1,1-dimethylpropyl group; C³-C⁶ cycloalkyl group such as cyclopropyl group, cyclobutyl group, cyclopentyl group, cyclohexyl group; a C³-C₆ cycloalkyloxy group such as cyclopropyloxy group, cyclobutyloxy group, cyclohexyloxy group, cyclohexyloxy group, such as methoxy

group, ethoxy group, propoxy group, isopropoxy group, butoxy group, isobutoxy group, tert-butoxy group, pentyloxy group, and isopentyloxy group; a C4-C7 cycloalkylalkoxyl group such as cyclopropylmethoxy group, cyclopentylmethoxy group; a C1-C5 alkylthic group such as methylthic group, ethylthic group, propylthio group, butylthio group, and pentylthio group; a C1-C5 alkylsulfonyl group such as methanesulfonyl group, ethanesulfonyl group, propanesulfonyl group, butanesulfonyl group, and pentanesulfonyl group; a halogen atom such as fluorine atom, chlorine atom, bromine atom, and iodine atom; a  $C_1$ - $C_5$  halogenated alkyl group such as trifluoromethyl group; hydroxyl group; cyano group; nitro group; formyl group; a C2-C5 alkylcarbonyl group such as acetyl group, propionyl group, butyryl group, and valeryl group; a benzene ring which may be substituted, a naphthalene ring which may be substituted, an optionally substituted heterocyclic ring having 1 to 4 hetero atoms selected from the group consisting of oxygen atom, sulfur atom and nitrogen atom, and having 5 to 10 ring-constituting atoms in total; a phenoxy group which may be substituted; a phenylamino group which may be substituted; an amino group; a C1-C5 monoalkylamino group such as methylamino group, ethylamino group, propylamino group, isopropylamino group, butylamino group, isobutylamino group, tert-butylamino group, pentylamino group, and isopentylamino group; a C2-C10 dialkylamino group such as dimethylamino group, ethylmethylamino group, diethylamino group, methylpropylamino group, and diisopropylamino group; a C1-C5 monoalkylaminomethyl group such as methylaminomethyl group, ethylaminomethyl group, propylaminomethyl group, isoproylaminomethyl group, butylaminomethyl group, isobutylaminomethyl group, tert-butylaminomethyl group, pentylaminomethyl group, isopentylaminomethyl; a  $C_2$ - $C_{10}$  dialkylaminomethyl group such as dimethylaminomethyl group, diethylaminomethyl group, ethylmethylaminomethyl group, methylpropylaminomethyl group; pyrrolidinylmethyl group; piperidinylmethyl group; morpholinomethyl group; piperazinylmethyl group; pyrrolylmethyl group; imidazolylmethyl group; pyrazolylmethyl group; and triazolylmethyl group.

When the C<sub>6</sub>-C<sub>10</sub> aryl ring represented by Y¹ has one or more substituents, the ring may be substituted by one or more substituents selected from the groups consisting of halogen atoms, a C₁-C₅ alkyl group, a C₃-C₆ cycloalkyl group, a C₃-C₆

cycloalkyloxy group, a C<sub>1</sub>-C<sub>5</sub> alkoxy group, a C<sub>4</sub>-C<sub>7</sub> cycloalkylalkoxy, a C<sub>1</sub>-C<sub>5</sub> alkylthio group, a C<sub>1</sub>-C<sub>5</sub> alkylsulfonyl group, a C<sub>1</sub>-C<sub>5</sub> halogenated alkyl, and a benzene ring.

When the ring represented by X, X1 or Y1 has one or more substituents, the substituent may further have one or more substituents selected from the group consisting of a C1-C5 alkyl group such as methyl group, ethyl group, propyl group, isopropyl group, butyl group, isobutyl group, sec-butyl group, tert-butyl group, pentyl group, isopentyl group, neopentyl group, 1,1-dimethylpropyl group; C3-C6 cycloalkyl group such as cyclopropyl group, cyclobutyl group, cyclopentyl group, cyclohexyl group; a C3-C6 cycloalkyloxy group such as cyclopropyloxy group, cyclobutyloxy group, cyclopentyloxy group, cyclohexyloxy group; hydroxy group; a C1-C5 alkoxy group such as methoxy group, ethoxy group, propoxy group, isopropoxy group, butoxy group, isobutoxy group, tert-butoxy group, pentyloxy group, and isopentyloxy group; a C4-C7 cycloalkylalkoxy group such as cyclopropylmethoxy group, cyclopentylmethoxy group; a C1-C5 alkylthio group such as methylthio group, ethylthio group, propylthio group, butylthio group, and pentylthio group; a C1-C5 alkylsulfonyl group such as methanesulfonyl group, ethanesulfonyl group, propanesulfonyl group, butanesulfonyl group, and pentanesulfonyl group; a halogen atom such as fluorine atom, chlorine atom, bromine atom, and iodine atom; a C1-C5 halogenated alkyl group such as trifluoromethyl group; a C1-C5 halogenated alkoxy group such as trifluoromethoxy group, 2,2,2-trifluoroethoxy group; hydroxyl group; cyano group; nitro group; formyl group; a C2-C6 alkylcarbonyl group such as acetyl group, propionyl group, butyryl group, and valeryl group; amino group; a C1-C5 monoalkylamino group such as methylamino group, ethylamino group, propylamino group, isopropylamino group, butylamino group, isobutylamino group, tert-butylamino group, pentylamino group, and isopentylamino group; a C2-C10 dialkylamino group such as dimethylamino group, ethylmethylamino group, diethylamino group, methylpropylamino group, and diisopropylamino group; a C2-C10 monoalkylaminomethyl group such as methylaminomethyl group, ethylaminomethyl group, propylaminomethyl group, isoproylaminomethyl group, butylaminomethyl group, isobutylaminomethyl group, tert-butylaminomethyl group, pentylaminomethyl group, isopentylaminomethyl; a C3-C11

dialkylaminomethyl group such as dimethylaminomethyl group, diethylaminomethyl group, ethylmethylaminomethyl group, methylpropylaminomethyl group and the like.

R may preferably be a C<sub>1</sub>-C<sub>8</sub> alkyl group, more preferably a methyl group or an ethyl group. The substituent of the alkyl group may preferably be a C<sub>3</sub>-C<sub>8</sub> alkyl group.

X may preferably be a benzene ring which may be substituted, a benzyl group which may be substituted, a naphthyl group which may be substituted, a benzofuran ring which may be substituted, a dihydrobenzofuran ring which may be substituted, a benzisoxazole ring which may be substituted, a benzisoxazole ring which may be substituted, a benzisothiazole ring which may be substituted, a benzisothiazole ring which may be substituted, and a benzopyrazole ring which may be substituted; more preferably a benzene ring which may be substituted, a benzyl group which may be substituted. Substituent of X may preferably be selected from the group consisting of a halogen atom, a C1-C4 alkyl group, a C1-C4 alkoxy group, a hydroxy group, a nitro group, a cyano group, a perhalogenated C1-C4 alkyl group, a carboxyl group, a C1-C4 alkoxycarbonyl group, a C1-C4 alkylthio group, a C1-C4 alkoxysulfonyl group, and amino group which may be substituted by a C1-C4 alkyl group.

The compounds represented by the aforementioned formula (I) may form a salt. Examples of the salt include, when an acidic group exists, salts of alkali metals and alkaline earth metals such as lithium, sodium, potassium, magnesium, and calcium; salts of ammonia and amines such as methylamine, dimethylamine, trimethylamine, dicyclohexylamine, tris(hydroxymethyl)aminomethane, N.N-bis(hydroxyethyl)piperazine, 2-amino-2-methyl-1-propanol, ethanolamine, N-methylglucamine, and L-glucamine; or salts with basic amino acids such as lysine,  $\delta$ -hydroxylysine, and arginine. When a basic group exists, examples include salts with mineral acids such as hydrochloric acid, hydrobromic acid, sulfuric acid, nitric acid, phosphoric acid; salts with organic acids such as methanesulfonic acid, benzenesulfonic acid, p-toluenesulfonic acid, acetic acid, propionic acid, tartaric acid, fumaric acid, maleic acid, malic acid, oxalic acid, succinic acid, citric acid, benzoic acid, mandelic acid, cinnamic acid, lactic acid,

glycolic acid, glucuronic acid, ascorbic acid, nicotinic acid, and salicylic acid; or salts with acidic amino acids such as aspartic acid, and glutamic acid.

In addition to the 3-substituted-4-pyrimidone derivatives represented by the aforementioned formula (I) and salts thereof, their solvates and hydrates also fall within the scope of the present invention. The 3-substituted-4-pyrimidone derivatives represented by the aforementioned formula (I) may have one or more asymmetric carbon atoms. As for the stereochemistry of such asymmetric carbon atoms, they may independently be in either (R) and (S) configuration, and the pyrimidone derivative may exist as stereoisomers such as optical isomers, or diastereoisomers. Any stereoisomers in a pure form, any mixtures of stereoisomers, racemates and the like fall within the scope of the present invention.

Preferred compounds of the present invention are represented by formula (II):

$$(X)_{p} \qquad (II)$$

$$(X)_{q} \qquad (Y)_{r}$$

wherein R, X, Y are the same as those defined above; p is 0 or 1; q is 0 or 1; r is an integer of 0 to 6; p+q is 1 or 2;

and Z represents N or CZ1 wherein Z1 represents hydrogen atom or Y.

Examples of more preferred classes of compounds represented by formula (II) include:

- (1) those wherein R represents a C<sub>1</sub>-C<sub>3</sub> alkyl group which may be substituted by a C<sub>3</sub>-C<sub>6</sub> cycloalkyl group;
- (2) the compounds of the above (1) wherein R is methyl group or ethyl group; Y is in 3-, 4- or 5-position of the piperazine ring or the piperidine ring; p+q is 1; and r is an integer of 0 to 3;
- (3) the compounds of the above (2) wherein X is a C<sub>1</sub>-C<sub>8</sub> alkyl group which may be substituted or a C<sub>6</sub>-C<sub>10</sub> aryl ring which may be substituted; Y is a C<sub>1</sub>-C<sub>6</sub> alkyl group which may be substituted; p is 1; q is 0; r is an integer of 0 to 3; and Z is N or CH;

- (4) the compounds of the above (3) wherein X is a benzene ring which may be substituted, a benzyl group which may be substituted; Y is a methyl group which may be substituted; Z is N and r is 0 or 1;
- (5) the compounds of the above (2) wherein X is a benzene ring which may be substituted, a benzyl group which may be substituted, a benzoyl group which may be substituted, or a benzisothiazol ring which may be substituted; Y is a methyl group which may be substituted; Z is N and p is 0;
- (6) the compounds of the above (2) wherein X is a C<sub>1</sub>-C<sub>8</sub> alkyl group substituted by a benzene ring which may be substituted or a benzene ring which may be substituted; Y is a hydroxy group, a cyano group, or Y¹-CO- wherein Y¹ is a C<sub>1</sub>-C<sub>8</sub> alkyl group; Z is CH or C-Y and r is 0 or 1; and
- (7) the compounds of the above (6) wherein X is a benzyl group which may be substituted or a benzene ring which may be substituted; Y is a hydroxy group, a cyano group, or an acetyl group; Z is CH or C-Y and r is 0 or 1.

Examples of particularly preferred classes of compounds represented by formula (II) include:

- (1) those wherein R is methyl group, Y is CH<sub>3</sub>O-CO- group or CH<sub>3</sub>CH<sub>2</sub>O-CO- group, Z is N, p is 0, q is 1, r is 0 or 1 and Y is in 3-position of the piperazine ring;
- (2) those wherein R is methyl group, Y is methyl group, benzyl group or acetyl group, Z is N, p is 1, q is 0, r is 0 or 1 and Y is in 4-position of the piperazine ring;
- (3) those wherein R is methyl group, Y is methyl group, Z is N, p is 1, q is 0, r is 1 to 3 and Y is in 3-, 4-, or 5-position of the piperazine ring;
- (4) those wherein R is methyl group, Y is hydroxyl group or cyano group, Z is CH, p is 1, q is 0, r is 0 or 1 and X and Y are attached on the same carbon atom;
- (5) those wherein R is methyl group, Y is hydroxyl group, cyano group or acetyl group, Z is C-Y, p is 0, q is 1 and r is 1.

Examples of preferred compounds of the present invention are shown in the tables below. However, the scope of the present invention is not limited to the following compounds.

able-1				**************************************	<del></del>	
		R <sup>2</sup> R <sup>2</sup> N N O R <sup>1</sup>				
No.	R1	R2	R3	R4	R5	R6
A1	CH3-	H	H	CH3-	H	H
A2 A3	CH3-	H	H	CH3CH2-	н н	н
A4	СН3-	н.	Ĥ	Y	н	н
A5	CH3-	н	Н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н
A6	снз-	Н	н	人人	н	н
A7	снз-	н	Н		н	н
A8	СН3-	Н	н		Н	н
A9	СН3-	H	Н		Н	H
A10	СН3-	Н	H .	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	Н
A11	CH3-	H	H	<u>大人</u>	H	Н
A12	CH3-	н	н	7.	н	Н
A13	CH3-	Н	Н		Н	н
A15	CH3-	н	н	~~~\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н
A16	CH3-	Н	н	Y~~~	н	н
A17	СН3-	Н	н	n-C8H17-	Н	н
A18	CH3-	Н	Н	L~~~	н	н
A19	СН3-	н	н		Н	н
A20	СН3-	н	н		н	н
A21	снз-	Н	н		Н	н
A22	CH3-	н	н	<b>&gt;</b> →	Н	н
A23	снз-	н	н	$\Diamond$	Н	н
A24	снз-	н	н	$\bigcirc$	н	Н
A25	снз-	н	н		н	н

No.	R1	R2	[R3_	R4	R5	R6
A26	СН3-	Н	н	OH:	н	н
A27	снз-	н	н	<u></u>	н	Н
A28	снз-	Н	н	<b>-</b>	н	н
A29	СН3-	н	н		н	н
A30	СН3	н	н	F-(-)-1	н	н
A31	снз-	н	н	CI →	н	н
A32	снз-	н	н	CI	н	н
A33	снз-	Н	н	c <del></del> (>-₁	Н	н
A34	СН3-	н	Н	Br	Н	н
A35	СН3-	Н	н	Br.	н	н
A36	снз-	Н	н	Br—{_}}	н	Н
A37	снз-	н	н	<b>-</b> 1	н	н
A38	CH3-	н	н	<b>├</b>	н	н
A39	CH3-	Н	Н		н	н
A40	CH3	н	Н	CH₃ CH₃	н	н
A41	CH3-	н	н		Н .	н
A42	СН3-	н .	н	H <sub>3</sub> C-{}	н	н
A43	СН3-	н	Н	C <sub>2</sub> H <sub>5</sub> —{	Н	н
A44	СН3-	н	н	n-C <sub>3</sub> H <sub>7</sub> -		Н
A45	СН3-	н	н	n-C₄H <sub>9</sub> {_}-{	н	н
A46	CH3-	Н	н	OH C	н	н
A47	СН3-	н	н	HO —	н	н

No.	R1	R2	R3	R4	R5	R6
140.	17.1	11/2			11.0	1,70
A48	CH3	Н	н	HO-{_}}-{	Н	н
A49	снз-	н	н	OCH <sub>3</sub>	Н	н
A50	снз-	н	н	H₃CO ——{	н	н
A51	снз-	Н	н	H₃CO <del>-</del> ⟨_ <del>}-</del>	Н	н
A52	снз-	н	Н	C <sub>2</sub> H <sub>5</sub> O-{}-{	Н	н
A53	снз-	н	н	n-C₃H <sub>7</sub> O- <b>⟨</b> _}-	Н	н
A54	снз-	н	Н	n-C <sub>4</sub> H <sub>9</sub> O-	Н	н
A55	снз-	н	н .	NO <sub>2</sub>	н	н
A56	CH3-	Н	н	O <sub>2</sub> N 	Н	н
A57	CH3-	Н	н	O <sub>2</sub> N-(	н	Н
A58	CH3-	Н	н	CN	Н	н
A59	СН3-	н	Н	NC	Н	н
A60	CH3-	н	н	NC-()	н	н
A61	CH3-	н	н	CF <sub>3</sub>	н	н
A62	CH3-	н	Н	F <sub>3</sub> C	н	н
A63	снз-	Н	н	F <sub>3</sub> C-{}-{	н	н
A64	снз-	н	н	COOH	н	Н
A65	СН3-	н .	Н	HOOC	н	н
A66	снз-	н	н	H00C-{\rightarrow}-{\rightarrow}-{\rightarrow}	н	н
A67	снз-	н	н	CO <sub>2</sub> Me	н	н
A68	снз-	Н	н	MeO <sub>2</sub> C	н	н
A69	СН3-	н .	н	MeO <sub>2</sub> C-{}	н	н

No.	R1	R2	R3	R4	R5	R6
A70	CH3-	н	н	CO <sub>2</sub> Et	Н	н
A71	СН3-	Н	н	EtO <sub>2</sub> C	н	н
A72	снз-	н	н	EtO <sub>2</sub> C-{}	Н	Н
A73	СН3-	н	н	SMe	н	Н
A74	СН3-	Н	н	MeS	н	Н
A75	снз-	н	н	MeS-{	Н	н
A76	СН3	Н	н	SO₂Me	н	н
A77	снз-	н	н	MeO <sub>2</sub> S	Н	н
A78	снз-	н	н	MeO₂S-{_}	н	н
A79	СН3-	н	н	NH <sub>2</sub>	н	н
A80	СН3-	н .	н	H <sub>2</sub> N	н .	н
A81	снз-	Н	н	H <sub>2</sub> N-\(\bigcirc\)	Н	н
A82	CH3-	Н	н	NMe <sub>2</sub>	н	н
A83	снз-	н	н	Me₂N	н	н
A84	снз-	н	н	Me <sub>2</sub> N-{	н	н
A85	снз-	н	. Н	$\Box$	н	н
A86	снз-	н	н	CC,	н	н
A87	СН3-	н	Н	HZ Ž	н	н
A88	СН3-	Н	Н	HN	Н	н
A89	СН3-	н	н	0,	н	н
A90	CH3-	н	н	62,	н	н
A91	снз-	Н	н	ST	н	H.

No.	R1	R2	R3	R4	R5	R6
A92	CH3-	Н	н	s Z	н	н
A93	CH3-	н	н	HNN	н	Н
A94	CH3-	Н	н	HN	н	Н
A95	снз-	н	Н	HN HN	н	Н
A96	снз-	н	н	(Z)	н	Н
A97	снз-	н	н	0N > y	н	Н
A98	снз-	н	Н	N-J	н	н
A99	CH3-	н	н	N-O	н	н
A100	снз-	н	н	SN	Н	н
A101	снз-	н	н	N= S	н	Н
A102	снз-	н	н	NS	н	н
A103	СН3-	н	н	Q,,	н	н
A104	снз-	н	н	(N)	н	н.
A105	СН3	н	н	N, O,	Н	н .
A106	снз-	н	н	SN	Н	н
A107	снз-	н	н	S S	н	н
A108	снз-	н	н	N S	н	н
A109	снз-	н	н	CN <sup>→</sup>	н	Н
A110	снз-	н	н	<b>N</b> →	н	н
A111	снз-	н	н	N_	н	Н
A112	снз-	н	н	NN−4	н	н
A113	снз-	н	Н	N_N	н	н

No.	R1	R2	R3	R4	R5	R6
A114	снз-	н	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н .	н
A115	CH3-	н	Н		н	н
A116	СН3-	н	н		н	н
A117	СН3-	н	Н		н	н
A118	снз~	н	н	TON TON	н	н
A119	СН3-	н .	н	,CQ	н	н
A120	СН3-	н	н	Çî,	н	н
A121	СН3-	н	н		н	н
A122	СН3-	Н	н		н	н
A123	снз-	Н	Н	Ö	Н	н
A124	СН3-	н	н	'CC;	н	Н
A125	снз-	н	Н	,CC	н	н
A126	СН3-	н	н	Ŷ.	Н	н
A127	СН3-	н	н	(I)	н	н.
A128	снз-	н	н	O.	Н	н
A129	снз-	н	н		Н	Н
A130	СН3-	Н	н	TOIS	н .	н
A131	снз-	н	н	, Ds	н	н
A132	СН3-	н	Н	Ĉ.	Н	н
A133	CH3-	н	н		н	н
A134	снз-	н	н	J.W	н	H
A135	снз-	н	н		н	н

No.	R1	R2	R3	R4	R5	R6
A136	снз-	н	н .	, Din	Н	Н
A137	CH3-	н	н	Ţì,	н	Н
A138	СН3-	н	н		н	Н
A139	СН3-	н .	н	N N	н	Н
A140	снз-	н	н		н	Н
A141	снз-	н	н	CTN-1	Н	н
A142	СН3-	н	н	N.	н	н
A143	снз-	н	н	'CI'	н	н
A144	CH3-	н	н	, (C)	н	н
A145	снз-	Н	н		н	н
A146	СН3-	Н	Н	()s <sup>N</sup> →	H ·	н
A147	CH3-	н	н	N S	н	н
A148	снз-	н	н	TOIS	Н	н
A149	CH3-	H ,	н	, CJ s	н	н
A150	снз-	Н	н	Ç s	н	н
A151	снз-	Н	н		Н	н
A152	снз-	Н	н	Ö	н	н
A153	снз-	Н	н	TOW .	н	н
A154	СН3-	Н	н		н .	н
A155	снз–	Н	Н	Ĉ.	н	Н
A156	снз-	Н	н		Н	н
A157	снз-	н	н	ČÇ <sub>N</sub>	Н	н

No.	R1	R2	R3	R4	R5	R6
A158	СН3-	н	н	CLIN	н .	Н
A159	СН3-	н	н	,CT)	н	н
A160	CH3-	Н	Н	ĨŢ <sup>S</sup> N	н	н
A161	снз-	Н	Н	O <sup>i</sup> ,	Н	н
A162	снз-	н	н	المراجع	н	н
A163	снз–	н	н	FO <sup>1</sup> ,	н	Н
A164	СН3-	Н	н	-Oir	Н	Н
A165	снз–	н	н	Cl o',	н	н
A166	СН3-	Н	н	CI	н	н
A167	снз-	н	Н		Н	н
A168	СН3-	Н	н	Br O	н	н
A169	СН3-	н	н .	Brody	н	н
A170	СН3-	Н	н		н	н
A171	снз-	Н	н	CHO	н	н
A172	СН3-	Н	н	H3C C	н	н
A173	СН3-	н	н		Н	H
A174	СН3-	н	н	CH30 0	н	н
A175	CH3-	н	н	H <sub>3</sub> CO	н	Н
A176	CH3-	н	н		н	Н
A177	снз-	Н	Н	100	Н	н
A178	CH3-	Н	Н	02N	н	н
A179	СН3-	Н	н	N P	н	н

No.	R1	R2	R3	R4	R5	R6
				EH O		Ţ
A180	СН3-	н ,	Н	O,	Н	Н
A181	СН3-	Н	н	HO	н	н
A182	СН3-	н	н	HO. S. P.	н	н
A183	СН3-	н .	н	NH 34	н	н
A184	СН3-	н	н	H <sub>2</sub> N	н	н
A185	CH3-	н	Н	L NO I'M	н	Н
A186	СН3~	н	Н	EN O	н	н
. A187	СН3-	н	н	NC CYLY	н	н
A188	снз-	н	Н	NC JI',	Н	н
A189	снз-	н	н		Н	н
A190	снз-	н	н		н	н
A191	снз-	н	Н	<u></u>	Н	н
A192	снз-	н	Н	<b>)</b>	н	н
A193	СН3-	Н	н	~\ <sup>1</sup> },	Н	н
A194	снз-	Н	н	\rangle, \ra	Н	Н
A195	СН3-	н	н	کیا <sub>ب</sub>	н	н
A196	снз-	н	н	~~~~	Н	н
A197	СН3-	Н	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н
A198	СН3-	Н	н .	~~ <sup>1</sup> ,	н	н
A199	снз-	н	н	~~~°;	н	н
A200	СН3-	н .	н	~~~ <sup>1</sup> ,	н	н
A201	СН3-	H.	н	√ <sup>1</sup> ,	н	н

No.	R1	R2	R3	R4	R5	R6
A202	CH3-	н	н		н	н
A203	CH3-	н	н		н	н
A204	СН3-	н	н		н	н
A205	снз-	H³CO_≻.	н	Н	н	н
A206	снз-	H³CO, Դ.	Н	СН3-	Н	н
A207	снз-	H³CO_5.	н	снзсн2-	н	Н
A208	снз-	H³CO, 5	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н
A209	снз-	H <sub>3</sub> CO ×	н	7	н	н
A210	СН3-	H <sub>3</sub> CO >	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н
A211	CH3-	H³CO_^≻	н	Li	н	н
A212	СН3-	H₃CO ≻	н	$\gamma$	н	н
A213	снз-	O H₃CO ≻	н	Y	н	н
A214	снз-	H <sub>3</sub> CO >	н	~~\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	Н
A215	СН3-	н₃со	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	Н
A216	СН3-	H₃CO ≻	н	人、	н	н
A217	СН3-	.H₃CO →	H.	7	н	н
A218	СН3-	H <sub>3</sub> CO >	н	~~~`	н	н
A219	снз-	H₃CO →	н		н	н
. A220	снз-	H³CO, ≻	н	~~~	入 H .	н
A221	сн3-	H <sub>3</sub> CO >	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	ч	н
A222	СН3-	H³CO, λ	н	n-C8H17-	н	н
A223	СН3-	H <sub>3</sub> CO '	н	J.	入 H	н

No.	R1	R2	R3	R4	R5	R6
A224	СН3-	H³CO_≻ O	н	Q	н	н
A225	0Н3-	H₃CO ≻	н		н	н
A226	СН3-	H³CO_≻	н		Н	н
A227	СН3-	H³CO_,	н		н	н
A228	снз-	H <sub>3</sub> CO '	н	$\Diamond$	н	н
A229	СН3-	H³CO >	н		Н	Н
A230	СН3~	H³CO,	н		н	н
A231	CH3-	H³CO →	н	$\bigcirc$	Н	н
A232	снз-	н₃сой	н	<b>◯</b> ─¹	н	н
A233	СН3-	н₃сой	н		н	н
A234	CH3~	H <sub>3</sub> CO ×	н	F	н	н
A235	СН3-	H3CO_>	н	F-(-)(	Н	н
A236	СН3-	H₃CO y	н	CI →	н	н
A237	СН3-	H³CO, ≻	н	CI	н	н
A238	СН3-	H <sub>3</sub> CO >	н	c⊢ <b>∕</b>	н	н
A239	СН3-	H³CO, >	н	Br	н	н
A240	СН3-	H³CO_^^	н	Br.	н	н
A241	СН3-	H <sub>3</sub> CO /	н	Br─⟨	Н	н
A242	СН3-	H <sub>3</sub> CO y	н	CH₃	Н	Н
A243	СН3-	H <sub>3</sub> CO >	н	H <sub>3</sub> C	н	н
A244	снз-	H <sub>3</sub> CO >	н	H <sub>3</sub> C-{_}-{	н	н
A245	онз-	H3CO, X	н	C <sub>2</sub> H <sub>5</sub> -{}-{	н	н

No.	RI	R2	R3	R4	R5	Re
A246	CH3-	H3CO_>	н	л-C <sub>3</sub> H <sub>7</sub> -⟨	н	н
A247	снз-	H3CO,	н	n-C <sub>4</sub> H <sub>9</sub> -	н	н
A248	CH3-	H <sub>3</sub> CO >	н	OCH₃	н	н
A249	снз-	H₃CO →	н	H₃CQ —}-{	н	н
A250	СН3-	H³CO_}.	н	H₃CO-{_}}-{	н	н
A251	CH3-	H³CO_>	н	C <sub>2</sub> H <sub>5</sub> O-{_}-{	Н	н
A252	снз-	H³CO_}\	н	n-C <sub>3</sub> H <sub>7</sub> O-	н	н
A253	СН3-	H <sub>3</sub> CO >	н	n-C <sub>4</sub> H <sub>9</sub> O-	н	н
A254	СН3-	H <sub>3</sub> CO 7	н	NO <sub>2</sub>	Н	н
A255	СН3-	H <sub>3</sub> CO >	Н	O <sub>2</sub> N	н	н
A256	CH3-	H³CO_}\	Н	O <sub>2</sub> N-(	Н	н
A257	снз-	H₃CO →	н	CN	н	н
A258	СН3-	H³CO,≻	н	NC	н	н
A259	снз-	H³CO, ≻	н :	ис-⟨∑─┤	н	н
A260	снз-	H₃CO 7	н	NMe <sub>2</sub>	н	н
A261	СН3-	H₃CO 7	Н	Me <sub>2</sub> N	Н	н
A262	снз-	H³CQ_>	н	Me <sub>2</sub> N-⟨¯¯	н	н
A263	снз-	H³CO_}	н		н	н
A264	СН3-	H₃CO y	н	CC,	н	н
A265	СН3-	H <sub>3</sub> CO 'y	н	O <sup>i</sup> ,	н	н
A266	снз-	H³CO ≻	н	Qi,	н	н.
A267	CH3-	H³CO_}	н	OO!	н	н

No.	R1	R2	R3	R4	R5	R6
A268	CH3-	H <sub>3</sub> CO y	н	۵,	н	н
A269	CH3-	H³CO_}\	н	, , , , , , , , , , , , , , , , , , ,	Н	н
A270	СН3-	C₂H₅O →	н	н	н	н
A271	CH3~	C₂H₅O →	н	СН3-	н	н
A272	СН3-	O C₂H₅O →	н	CH3CH2~	н	н
A273	GH3-	O C₂H₅O →	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н
A274	CH3	C <sub>2</sub> H <sub>5</sub> O ,	н	$\nearrow$	Н	н
A275	CH3-	C <sub>2</sub> H <sub>5</sub> O, 7	Н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н
A276	СН3-	O C₂H₅O '-'	н	\\ \	Н	н
A277	СН3-	C <sub>2</sub> H <sub>5</sub> O	н	~~	Н	н
A278	CH3-	C <sub>2</sub> H <sub>5</sub> O ,	н	丫	н	Н
A279	CH3-	C <sub>2</sub> H <sub>5</sub> O	н		н	Н
A280	СН3-	C₂H₅O Y	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	н
A281	CH3-	C₂H₅O →	н	X.	н	н
A282	снз-	C₂H₅O →	н.	7	н	н
A283	СН3-	C₂H₅O →	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	н
A284	СН3-	C <sub>2</sub> H <sub>5</sub> O 7	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н
A285	снз-	C₂H₅O →	н	~~~`\	н	н
A286	снз-	C₂H₅O r	н	Y~~~	н .	н
A287	CH3-	C₂H₅O →	н	n-C8H17-	н	н
A288	СН3-	C <sub>2</sub> H <sub>5</sub> O '' '' C <sub>2</sub> H <sub>5</sub> O '' '' ''	н		н	н
A289	СН3-	C₂H₅O →	н		н	н

No.	R1	R2	R3	R4	R5	R6 ·
A290	снз-	C <sub>2</sub> H <sub>5</sub> O >	н		н	н
A291	снз-	C₂H₅O Y	н		Н	н
A292	снз-	C <sub>2</sub> H <sub>5</sub> O y	н	<b>⊳</b> }	Н	н
A293	CH3-	C₂H₅O y	н	$\Diamond$	Н	н
A294	снз-	C₂H₅O <sup>™</sup> ≻∕	н	$\bigcirc$	н ,	н
A295	снз-	C <sub>2</sub> H <sub>5</sub> O y	н		н	н
A296	СН3-	C <sub>2</sub> H <sub>5</sub> O ,	н		Н	н
A297	снз-	C <sub>2</sub> H <sub>5</sub> O ' >	н		Н	н
A298	СН3-	C₂H₅O →	н	<b>_</b> -	Н	н
A299	СН3-	C <sub>2</sub> H <sub>5</sub> O y	н		Н	н
A300	снз-	C <sup>2</sup> H <sup>2</sup> O >	н		Н	н
A301	снз-	C <sub>2</sub> H <sub>5</sub> O <sup>11</sup> >	н	CI	Н	н
A302	снз-	C <sub>2</sub> H <sub>5</sub> O	н	CI	Н	Н
A303	СН3-	O C₂H₅O →	н	CI-{}-{	н .	н
A304	снз-	C₂H₅O →	н	Br	Н	Н
A305	снз-	C <sub>2</sub> H <sub>5</sub> O -	Н	Br	Н	Н
A306	CH3-	C <sub>2</sub> H <sub>5</sub> O ,	н	Br—{}	Н	Н
A307	СН3-	C₂H₅O Y	н	CH <sub>3</sub>	Н	Н
A308	СН3-	C₂H <sub>6</sub> O y	н	H <sub>3</sub> C ————————————————————————————————————	Н	Н
A309	СН3-	C <sub>2</sub> H <sub>5</sub> O >	н.	H <sub>3</sub> C:-{_}}{	Н	Н
A310	снз-	C <sub>2</sub> H <sub>5</sub> O 7	н	C <sub>2</sub> H <sub>5</sub> -{}-{	н	н
A311	снз-	C <sub>2</sub> H <sub>5</sub> O 7	н	n-C <sub>3</sub> H <sub>7</sub> {}{	н	н

No.	R1	R2	IR3	R4	R5	R6
A312	CH3-	C <sub>2</sub> H <sub>5</sub> O /	н	n-C <sub>4</sub> H <sub>9</sub> {}-{	н	н
A313	СН3-	C₂H₅O →	н	OCH₃	н	н
A314	СН3-	C <sub>2</sub> H <sub>5</sub> O 7	н	H <sub>3</sub> CQ	н	н
A315	снз-	C₂H₅O Y	н	H <sub>3</sub> CO-{_}-{	н	Н
A316	снз-	O C₂H₅O →	н	C <sub>2</sub> H <sub>5</sub> O-	н	н
A317	снз-	C₂H₅O →	н	n-C₃H <sub>7</sub> O-⟨}{	н	н
· A318	снз-	C <sub>2</sub> H <sub>5</sub> O /	н	n-C <sub>4</sub> H <sub>9</sub> O-{}-;	н	H
A319	снз-	C₂H₅O →	н	NO <sub>2</sub>	н	н
A320	снз–	C <sub>2</sub> H <sub>5</sub> O >	Н	O <sub>2</sub> N	н	Н
A321	СН3-	C₂H₅O	Н		н	н
A322	снз-	C₂H₅O ×	н	CN CN	Н	н
A323	СН3-	O C₂H₅O →	Н	NC	н	н
A324	СН3-	O C₂H₅O ≻	Н	NC-	н	н
A325	снз-	C <sub>2</sub> H <sub>5</sub> O →	н	NMe <sub>2</sub>	н	Н
A326	СН3-	C₂H₅O →	Н	Me <sub>2</sub> N	н	Н
A327	снз-	C <sub>2</sub> H <sub>5</sub> O Y	н .	Me <sub>2</sub> N-	н	н
A328	снз-	C <sub>2</sub> H <sub>5</sub> O →	н		н	Н
A329	снз-	C₂H <sub>6</sub> O →	Н	CC,	Н	н
A330	снз-	C <sub>2</sub> H <sub>5</sub> O >	н	O <sup>i</sup>	н	н
A331	СН3-	C <sub>2</sub> H <sub>5</sub> O <sup>T</sup> >	н	Qi,	н	н .
A332	GH3~	O C <sub>2</sub> H <sub>5</sub> O 7	H	OOL	н	н
A333	снз-	C <sub>2</sub> H <sub>5</sub> O ×	н	2,	н	н

No.	R1	R2	R3	R4	R5	R6
A334	снз-	C <sub>2</sub> H <sub>5</sub> O 7	Н	<u>,</u>	н	н
A335	CH3-	,	H	н	н	н
A336	CH3-	CH3CH2-	Н	н	Н	н
A337	СН3-	<b>/</b> \\	Н	н	н	н
A338	CH3-	$\rightarrow$	н	н	Н	н
A339	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	Н	н	н
A340	снз-	L	Н	Н	Н	н
A341	СН3-	~	н	Н	Н	н
A342	СН3-	Y	н	н .	н	Н
A343	СН3-	~~\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	Н .	Н	Н
A344	Снз-	Y~~`	н	Н	н	н
A345	СН3-	人人	н	Н	н	Н
A346	CH3-	7	н	Н	Н	н
A347	CH3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	н	Н
A348	снз-		н	н	н	Н
A349	снз-	~~~``	H	н	н	н
A350	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	Н	Н
A351	CH3-	n-C8H17-	н	н	н	н
A352	СН3-	L	Н	н	н	н
A353	СН3-		Н	н	н	н
A354	СН3-		Н	н	н	н
A355	снз-	Qui	н	Н	н	Н

No.	R1	R2	R3	R4	R5	R6
A356	СН3-	$\triangleright$	Н	Н		н
A357	СН3-	<u> </u>	н	Н	Н	H
A358	СН3-	$\bigcirc \dashv$	Н	н	н	н
A359	СН3-	$\bigcirc$ $\dashv$	н	н	н	н
A360	снз-	$\bigcirc$ -i	н	н	н	н
A361	СН3-	<u></u>	н	Н	Н	Н
A362	CH3~		Н	Н	Н	н
A363	CH3-	<u></u>	н	н	н	н
A364	снз-	<b>-</b>	н	н	н	Н
A365	CH3-	F	н	н	н	н
A366	СН3-	F-();	н	н	Н	н
A367	СН3-	F-(>-1	Н	н	н	н
A368	CH3-	F——	н	Н	Н	н
A369	снз-	CI →	н	Н	н	н
A370	онз-	CI	Н	н	н	н
A371	снз-	c <del>(-)-</del> -	н	н	н	Н
A372	онз-	CH_	н	Н	н	н
A373	CH3	CH	н	н	н	н
A374	снз-	Br	н	н	н	н
A375	СН3-	Br.	н	н	Н	н
A376	СН3-	Вг−СУ−{	н	н	н	Н
A377	CH3	Br{}	н	н	н	н

No.	R1	R2	R3	R4	R5	R6
A378	СН3-	Br—{	н	н	н	н
A379	CH3~		Н	н	Н	н
A380	снз-		Н	н	Н	н
A381	снз-	<del></del>	н	Н	н	Н
A382	снз-	(/;	I	Н	н	H
A383	СН3-	H <sub>3</sub> C	н	н	Н	н
A384	снз-	H <sub>3</sub> C-{	н	н	н	н
A385	CH3-	C <sub>2</sub> H <sub>5</sub> —{	н	н	Н	н
A386	СН3-	n-C <sub>3</sub> H <sub>7</sub> -	Н	Н	Н	н
A387	снз-	n-C <sub>4</sub> H <sub>9</sub> -	H	н	Н	н
A388	снз-	OH OH	н	н	н	Н
A389	СН3-	HO	Н	н	Н	Н
A390	снз-	HO-{\bar{\bar{\bar{\bar{\bar{\bar{\bar	н	н	н	н
A391	CH3-	OCH <sub>3</sub>	Н	Н	н	Н
A392	снз-	H <sub>3</sub> CO	н	н	н	н
A393	снз-	H₃CO-⟨}-{	н	н .	н	н
A394	снз-	H₃CO- <b>(_)</b> {	н	н	Н	н
A395	CH3-	H <sub>3</sub> CO-()III-\$	Н	н	н	н
A396	CH3-	OC <sub>2</sub> H <sub>5</sub>	н .	Н	Н	н
A397	СН3-	C <sub>2</sub> H <sub>5</sub> Q	Н	н	Н	н .
A398	CH3-	C <sub>2</sub> H <sub>5</sub> O-	н	Н	н	н
A399	CH3-	n-C <sub>3</sub> H <sub>7</sub> O-⟨	н	Н	н	н

No.	RI	R2	R3	R4	R5	R6
A400	CH3-	n-C <sub>4</sub> H <sub>9</sub> O-	н	н	н	Н
A401	снз-	NO <sub>2</sub>	н	н	Н	н
A402	СН3-	O <sub>2</sub> N	н	н	н	н
A403	СН3-	O <sub>2</sub> N-(	н	н	н	н
A404	CH3-	CN C	н	н	н	н
A405	СН3-	NC	н	н	н	Н
A406	СН3-	NC-{}-	н	н	н	н
A407	СН3-	CF <sub>3</sub>	Н	Н	н	н
A408	снз-	F <sub>3</sub> C	н	Н	н	н
A409	снз-	F <sub>3</sub> C-{_}-{	н	Н	н .	н
A410	СН3-	COOH COOH	н	н	Н	н
A411	СН3~	HOOC.	н	н	н	н
A412	снз-		н	Н	н	н
A413	снз-	CO₂Me	н	н	н	н
A414	снз-	MeO <sub>2</sub> C	н	Н	н	н
A415	СН3	1	Н	н	н	н
A416	CH3		н	н	н	н
A417	снз-	EtO <sub>2</sub> C	Н	н	н .	н
A418	СН3-	1	н	н	н	н
A419	снз-		н	н	н	Н
A420	снз-	MeS	н	Н	Н	Н
A421	снз-	MeS-{_}-{	Н	Н	н	н

No.	R1	R2	R3	R4	R5	R6
A422	CH3-	SO₂Me	н	н	н	н
A423	СН3-	MeO <sub>2</sub> S	н	н	Н	н
A424	CH3-	MeO <sub>2</sub> S-{_}-{	H	н	н	н
A425	снз-	NH <sub>2</sub>	Н	н .	Н	Н
A426	CH3-	H <sub>2</sub> N	Н	Н	Н	н
A427	CH3-	H <sub>2</sub> N-{	н	н	н	Н
A428	снз-	NMe₂ →	н	н	Н	Н
A429	снз-	Me <sub>2</sub> N	н	Н	н	н
Ą430	СН3-	Me₂N-⟨¯¯⟩{	н	Н	н	н
A431	СН3-		Н	Н	н	Н
A432	снз-	(N-Q)	н	н	н	Н
A433	СН3-	(N-(_)-1	н	н	н	н
A434	СН3-		н	н	н	н
A435	СН3	O-Q	н	н	н	н
A436	СН3-	\n-\_\-1	н	н	н	н
A437	СН3-	<b>○</b> \	H.	н	н	н
A438	снз-	<b>○</b> N- <b>○</b>	н	н	н	н
A439	СН3-		н	н	н	н
A440	CH3-	H <sub>3</sub> CN N-	н	н	н	н
A441	CH3~	H <sub>3</sub> CN N-	Н	Н	н	н
A442	СН3-	H3CN_N-{}-{	н	н	н	н
A443	снз-	H <sub>3</sub> C CH <sub>3</sub>	н	н	н	Н

No.	R1	R2	R3	R4	R5	lne l
INO.	141	CH <sub>3</sub>	173	π4	หอ	R6
A444	СН3-	H <sub>3</sub> C-{}	н	Н	Н	н
A445	снз-	H³C	н	н	н	н
A446	СН3-	CH,	н	н	н	н
A447	СН3	1,30	н	Н	Н	н
A448	CH3-	H <sub>3</sub> C	н	Н	Н	н
A449	СН3-	<b>-</b>	Н	н	Н	н
A450	СН3-	F-{=}-}	н	н	н .	н
A451	СН3-	\$	Н	н	н	Н
A452	СН3-	¢,	Н	н	н	н
A453	CH3-	F-\	Н	Н	н	Н
A454	СН3-	F	H ·	н	Н	н
A455	СН3-	CI_CI	н	н	н	Н
A456	СН3-	α-√_}-;	Н	Н	н	н
A457	СН3-		Н	H,	н	Н
A458	СН3-	ci G	н	н	н	Н
A459	снз-	G, CI—⟨□}—;	Н	Н	н	н
A460	СН3-	a'	н	н	н	Н
A461	снз–	· • ·	н	н	H .	н
A462	СН3	300	Н	н	н	н
A463	СН3-	H²CQ	н	н	н	Н
A464	CH3-	, och <sup>3</sup>	Н	H .	Н	н
A465	СН3-	H₃CO-{}}	н	Н	Н	н

No.	R1	R2 H₃GQ	R3	R4	R5	R6
A466	снз-	H <sub>3</sub> CO	н	н	. н	н .
A467	снз-	F_OCH <sub>3</sub>	н	н	н	Н
A468	СН3-	OCH₃ F—⟨S)→	н	н	н	н
A469	снз-	OCH <sub>3</sub>	н	н	Н	Н
A470	снз-	OCH <sub>3</sub>	н	н	н	н
A471	СН3-	OCH <sub>3</sub>	н	Н	н	н
A472	снз-	Poch?	н	н	н	Н
A473	снз-	H <sub>3</sub> CO	н	Н	н	н
A474	снз-	H <sub>3</sub> CQ	н	Н	Н	н
A475	снз-	H <sub>3</sub> CO_F	Н	н	. н	н
A476	СН3-	H₃CO-⟨S	н	н	н	н
A477	снз-	H <sub>3</sub> CO	н	н	н	н
A478	СН3-	H₃CO-⟨S	н	н	н	н
A479	CH3-	CI_OCH₃	н	н	Н	н
A480	CH3-	CI—⟨□>→{	н	н	Н	н
A481	CH3-	GCH₃	Н	н	н	н
A482	CH3-	CI CI	н	н	н	н
A483	снз-	H³CO CI—⟨□}—	Н	н	н .	н
A484	CH3-	Ci Ci	н	н	н	Н
A485	GH3-	H <sub>3</sub> CO_CI	н	н	н	н
A486	СН3-	H₃CO-{\(\)}	н	н	Н	н
A487	CH3-	H <sub>3</sub> CO -	н .	н	н	н

No.	R1	R2	R3	R4	R5	R6
A488	снз-	CI H₃CO-⟨¯_}~{	н	Н	н	н
A489	снз-	F_CH <sub>3</sub> ·	н	н	н	н
A490	снз-	CH <sub>3</sub> F—⟨ →	н	Н	н	н
A491	снз-	CH;	н	Н	н	н
A492	снз-	CH,	н	н	н	н
A493	снз-	H <sub>3</sub> C F-(	н	н	н	н
A494	снз-	H <sub>2</sub> C	н	н	н	Н
A495	СН3-	H <sub>3</sub> C_F	н	н	н	Н
A496	снз-	H₃C-⟨□}	н	н	н	н
A497	снз-	H <sub>5</sub> C	н	н	н	н
A498	снз-	H <sub>3</sub> C-\\	н	н	н	Н
A499	СН3-	Br. OCH₃	н	н	Н	н
A500	СН3-	OCH₃ Br-√→	н	н	Н	н
A501	СН3-	OCH <sub>3</sub>	н	н	н	н
A502	СН3-	OCH <sub>2</sub>	н	Н	н	н
A503	снз-	H <sub>3</sub> CO Br—{}	н	Н	н	н
A504	СН3-	H <sub>3</sub> CO	. н	Н	н	н
A505	СН3-	H <sub>3</sub> CO_Br	н	н	н	н
A506	СН3-	H₃CO-€S-	н	н	н	н
A507	снз-	H₃CO Sr	Н	н	н	н
A508	СН3-	H₃CO-⟨	Н	н	Н	Н
A509	СН3-	H <sub>3</sub> CO	Н	н	н	н

No.	RI	R2	R3	R4	R5	R6
A510	CH3-	OCH <sub>3</sub>	н	Н	Н	н
A511	CH3-	CN-CD-OCH3	н	н	Н	н
A512	CH3-	H <sub>3</sub> CO >-	н	Н	н	н
A513	СН3	H <sub>3</sub> CO	н	н	н	н
A514	CH3-	Chry	н	н	н	Н
A515	СН3~		н	н	Н	н
A516	снз-	FC∑+,	н	н	Н	н
A517	СН3-	H³CO-{\$\sum_1\}	н	н	н	н
A518	СН3-	OCH <sub>3</sub>	н	н	н	н
A519	СН3-	OCH <sup>3</sup> H³CO-{∑}-1 OCH <sup>3</sup>	н	н	н	н
A520	СН3-	a-<\$\frac{\circ}{\circ};	Н	н	н	Н
A521	СН3-	CI-√∑H CI-√∑H	н	н	н	н
A522	снз-	H²co-{\$\frac{1}{2}}	н	н	н	н
A523	снз-	CI-(2)-5 OCH3	н	Н	Н	н
A524	СН3-	H³CO-⟨∑-1 OCH²	Н	Н	Н	Н
A525	СН3-	OCH <sub>3</sub>	Н	н	н	н
A526	снз-	H <sub>3</sub> CO	Н	Н	Н	н
A527	снз-	H <sub>3</sub> CO-{_}-{_}-{_}-{_}-{_}{_}{_}{_}{_	Н	н .	н	Н
A528	снз-	OCH <sub>3</sub> >	н	н	Н	н
A529	снз-	H <sub>3</sub> CO ,	н	Н	н	н
A530	СН3-	н₃со-⟨>-⟨>	н .	н	н	Н
A531	снз-	Q.Q	н	н	н	Н

No.	R1	R2	R3	R4	R5	R6
140.	,,,	H-co		DVT	110	110
A532	CH3-		н	Н	Н	н
A533	СН3-	H₃∞-	н	Н	н	н
A534	онз-	<b>₫-</b> 0-₁	Н	Н	н	н
A535	снз-	<b>E</b>	H	н	н	н
A536	снз-	F-(>-(>-1	н	н -	н .	н
A537	СН3-	ď-ď`	H	Н	Н	н
A538	СН3-		Н	н	н .	Н
A539	CH3-		н	Н	н	н
A540	СН3-	Q-6	Н	н	н	н
A541	СН3-	\$\dag{\omega}	Н	Н	н	н
A542	CH3		Н	H .	н	н
A543	CH3-		Н	Н	н	Н
A544	CH3-	CCT'	н	Н	н	н
A545	CH3-		н	н	н	н
A546	снз-	HNJ	Н	Н	н	н
A547	CH3-	S.	Н	н	н	н
A548	СН3-	d.),	н	н	н	н
A549	CH3-	St	н	н	н	н
A550	СН3-	s.J.,	н	н	Н	н
A551	снз-	HNN	н	н	Н	н
A552	снз–	HN	н	н	Н	н
A553	СН3	HN A	н	н	н	н

No.	R1	R2	R3	R4	R5	R6
A554	CH3-	N N	н	н	н	Н
A555	CH3-	and,	н	н	Н	н
A556	CH3-	N=)	н	Н	н	H
A557	CH3-		Н	н	н	н
A558	CH3-	S <sub>N</sub>	н	н	н	Н
A559	снз-	N= S	н	н	н	н
A560	снз-	N-S	Н	н	н	н
A561	снз-	√=N O, ✓, γ	н	н	н	н
A562	СН3-	O T	н	н	н	н
A563	СН3-	N Z	н	Н	Н	н
A564	CH3-	S-N	н	Н	н	H
A565	СН3-	S	н	Н	Н	н
A566	снз-	N T	Н	Н	н	н
A567	СН3-	C <sub>N</sub> →	н	н	н	н
A568	CH3-		н	Н	н	н
A569	СН3		н	Н	н	н
A570	снз-	N →	н	Н	н	н
A571	СН3-		н	н	н	н
A572	CH3-	N N	н	н	н	н
A573	СН3-		н	н	н	н
A574	снз-	CT)	н	н	Н	н .
A575	СН3-		н	Н	н	н

No.	R1	R2	R3	R4	R5	R6
140	17.	7. H	110		110	1.0
A576	CH3-		н	Н	Н	Н
A577	СН3-	,CT	н	н :	н 	н
A578	СН3-		н	н	н	н
A579	CH3-		н	н	н	н
A580	CH3-		н	н	н	н
A581	CH3		н	Н	н	н
A582	CH3~	TOO?	н	н	н	н
A583	CH3-	,CO	н	н	н	н
A584	CH3-	Ğ.	н	н	Н	н
A585	CH3-		н	н	Н	τ
A586	CH3-	CI)	н	н	н	н
A587	СН3	Ī;	н	Н	н	Н
A588	СН3-	TOIS	н	н	н	н
A589	CH3-	, Os	н	Н	н	н
A590	СН3-	Ĉ.	н	Н	н	Н
A591	снз-	CT.	н	н	н	н
- A592	снз-		н	Н	н .	н
A593	СН3-	TON	н	Н	Н	н
A594	снз-	'CLI'	Н	Н	Н	н
A595	СН3-	Ţ'n.	н	Н	Н	Н
A596	СН3-	(CTN) NH	Н	н	н	Н
A597	СН3-	Č',	н	н	н	Н

No.	R1	R2	R3	R4	R5	R6
140,	-	T. J.N	1.0		110	
A598	CH3-		Н	Н	н	н
A599	CH3-		н	н	н	н
A600	СН3-	<u> </u>	н	н	н	н
A601	CH3-	'CT'	H	н	н	н
A602	снз-	, Or	H	н	н	н
A603	СН3-	<b>P</b> 3	н	Н	н	н
A604	снз-	OTS+	Н	н .	н	н
A605	СН3-	Č,	н	Н	н	н
A606	СН3-	TO N	н	н	н	н
A607	CH3-	, O'S	н	н	н	н
A608	СН3-	Ž,	н	н	Н	н
A609	СН3~		н	Н	Н	н
A610	СН3-		н	н	н	н
A611	СН3-	TO31	н	Н	н	Н
A612	СН3-	,CT	н	н	н	н
A613	снз-	ĞÇ,	н	н	н	н
A614	снз-		н	н	н	н
A615	СН3-	ÖÇ,	н	н	Н	Н
A616	СН3	CA	н	н	н	н
A617	снз-	, CIÈN	Н	Н	н	Н
A618	снз-	ÇÇ"	н	н	н	н
A619	снз-	Çr.	н	н	н	Н

No.	R1	R2	R3	R4	R5	R6
		<b>~</b>				1
A620	CH3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	H	H	H	Н
A cn1	CH2-	·/^	ы	Li		,
A621	CH3-		н .	H	H	н
A622	CH3-	<u>~</u>	н	н	н	н
AUZZ	OFIS		11		n	n
A623	СН3-	CH3-	H	СНЗ	н	н
A624	CH3-	CH3CH2-	н	снз	н	н
		<u> </u>				
A625	CH3-	~~\ <sup>(</sup>	н	снз	н	н
		<u> </u>				
A626	CH3-	$\uparrow$	н	СНЗ	н	н
		·				
A627	СН3-	~ ~ `	н	снз	н	н
		1				
A628	CH3-	人入	н	СНЗ	н	н
		$\sim$				
A629	CH3-		Н	СНЗ	Н	Н
		<b>√</b> √				
A630	CH3~	1	н	СНЗ	H	Н
		^^\\				
A631	СН3-		Н	СНЗ	н	Н
4.633	CH3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	OUE	Н	
A632	UH3-			СНЗ		H
A633	CH3-	J z	Н	СНЗ	Н	н
			<u> </u>	0,10		
A634	снз-	$\sim$	н	СНЗ	H	н
					ļ	
A635	СН3-	<b>~~~</b> `.	н	СНЗ	н	н
	ļ	1				
A636	снз-	人~~	н	СНЗ	н	Н
		^ ^ 3	<b> </b>			ļ
A637	снз-		н	СНЗ	н	н
<b></b>	<del>                                     </del>	~~~\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\				<del> </del>
. A638	снз-		н	СНЗ	Н	н
A639	снз-	n-C8H17-	н	СНЗ	Н	Н
			1			1.
A640	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	H	СНЗ	H	Н
4044	0110		1.	O. I.S.		Ī.,
A641	CH3-		н	СНЗ	Н	Н

No.	R1	R2	R3	R4	R5	R6
A642	CH3-		Н	CH3	н	н
A643	CH3-		н	СН3	н	н
A644	снз-	D <del></del>	н	СНЗ	н	н
A645	CH3-	$\Diamond$ -1	н	снз	н	н
A646	снз-	$\bigcirc \dashv$	н	снз	Н	н
A647	CH3-	$\bigcirc$	н	снз	Н	н
A648	снз-	$\bigcirc$	н	СНЗ	н	н
A649	СН3-	<b>△</b> -₹	н	снз	Н	н
A650	СН3		н	снз	Н	Н
A651	снз-		н	СНЗ	Н	н
A652	СН3-	<u></u>	н	СНЗ	Н	н
A653	CH3~	F	н	СНЗ	Н	н
A654	СН3-	F-(	н	СНЗ	н	н
A655	CH3-		н	СН3	н	н
A658	снз-	F-\_\_\	н	СНЗ	н	н
A657	CH3-	CI	н	СН3	н	Н
A658	CH3-	CI.	н	СНЗ	н	н
A659	СН3-	c-(	н	снз	Н	н
A660	снз-	C-(-)-1	н	СНЗ	Н	н
A661	снз-	CI—(}III-{	н	СНЗ	н	н
A662	CH3-	Br	н	СНЗ	н	н
A663	CH3-	Br	н	снз	н	н

No.	R1	R2	R3	R4	R5	R6
A664	CH3~	Br—{}	н	СНЗ	н	н
A665	СН3-	Br-{}	Н	снз	н	н
A666	снз-	Br—⟨\in-{	н	снз	Н	н
A667	снз- :	<b>△</b> -₁	н	снз	н	Н
A668	снз-		н	СНЗ	н	н
A669	СН3-	<del></del>	н	снз	Н	Н
A670	СН3-	CH₃	н	СНЗ	Н	Н
A671	СН3-	H <sub>3</sub> C	Н	СНЗ	н	н
A672	снз-	H <sub>3</sub> C-{}-{	н	СНЗ	н	н
A673	CH3-	C <sub>2</sub> H <sub>5</sub> -{	Н	СНЗ	Н	н
A674	CH3-	n-C <sub>3</sub> H <sub>7</sub> -{{}}{	Н	СНЗ	н	н
A675	снз-	n-C <sub>4</sub> H <sub>9</sub> {	н	снз	н	н
A676	СН3-	OH .	Н	снз	н	н
A677	СН3-	HO	н	снз	н	н
A678	CH3-	HO-{\}-{	н	СНЗ	н	н
A679	СН3-	OCH₃	н	снз	н	н
A680	CH3-	H₃CQ	н	снз	Н	н
A681	CH3-	H₃CO- <b>⟨</b> _}~{	Н	снз	Н	н
A682	снз-	H <sub>3</sub> CO-{_}-{	н	снз	Н	н
A683	снз-	H <sub>3</sub> CO-	Н	снз	н	н
A684	снз-	OC <sub>2</sub> H <sub>5</sub>	н	СНЗ	Н	н
A685	снз-		н	CH3	н	н

No.	R1	R2	R3	R4	R5	R6
A686	СН3-		Н	снз .	н	н
A687	снз-	n-C <sub>3</sub> H <sub>7</sub> O-{}-{	н	снз	Н	н
A688	СН3-	n-C <sub>4</sub> H <sub>9</sub> O-⟨}-{	н	снз	н	н
A689	снз-	NO <sub>2</sub>	Н	снз	н	н
A690	СН3-	O <sub>2</sub> N	н	снз	н	н
A691	снз-	O <sub>2</sub> N-{}	н	снз	н	н
A692	СН3-	CN →	н	снз	Н	н
A693	СН3-	NC	Н	снз	н	н
A694	CH3-	NC-{}-{	н.	снз	н	н
A695	CH3-	CF <sub>3</sub>	н	снз	н	н
A696	снз-	F₃C ⟨_)(	н	снз	Н	н
A697	СН3-	F <sub>3</sub> C-{}-{	Н	снз	н	н
A698	CH3-	COOH	н	снз	Н	н
A699	CH3-	HOOC	н	снз	н	н
A700	CH3-	ноос-{}-	н	снз	н	Н
A701	CH3-	CO <sub>2</sub> Me	н	снз	н	н
A702	CH3-	MeO₂C	Н	СНЗ	н	н
A703	CH3-	MeO <sub>2</sub> C-{{}}	Н	снз	н	н
A704	CH3-	CO₂Et	н	снз	н	Н
A705	CH3-	EtO <sub>2</sub> C	н	СНЗ	н	Н
A706	снз-	EtO <sub>2</sub> C-{}-{	н	снз	н	н
A707	снз-	SMe	Н	снз	н	н

No.	R1	R2	R3	R4	R5	R6
A708	СН3-	MeS	н	СНЗ	Н	н
A709	СН3-	MeS-{_}-{	Н	CH3	н .	н
A710	СН3-	SO₂Me	н	СНЗ	Н	Н
A711	снз-	MeO <sub>2</sub> S	н	СНЗ	н	н
A712	снз-	MeO <sub>2</sub> S-	н	СНЗ	н	Н
A713	снз-	NH₂ →	н	снз	н	н
A714	снз-	H <sub>2</sub> N ⟨□}→	н	снз	н	Н
A715	снз-	H <sub>2</sub> N-\\\\\\\\	н	СНЗ .	Н	н
A716	СНЗ-	NMe <sub>2</sub>	н	СНЗ	Н	н
A717	снз-	Me <sub>2</sub> N	н	СНЗ	н	н
A718	снз-	Me <sub>2</sub> N-{	н	СНЗ	Н	н
A719	CH3-	(N-\)	н	СНЗ	н	Н
A720	СН3-	(N-Q)	Н	снз	н	н
A721	снз-	(n-()-i	Н	СНЗ	Н	н
A722	CH3-		н .	СНЗ	н	н
A723	CH3		Н	СНЗ	н	н
A724	СН3-		н	СНЗ	Н	н
A725	снз-		н	СНЗ	н	Н
A726	CH3-		н	СНЗ	н	н
A727	снз-	<b>○</b> N- <b>○</b> -1	н	СНЗ	H .	н
A728	CH3-	H3CN N-	Н	СНЗ	Н	н
A729	снз-	H3CN N-	н	СНЗ	н	н

No.	R1	R2	R3	R4	R5	R6
A730	CH3-	H3CN N-{}-{	н	CH3	н	н
A731	снз-	H <sub>3</sub> C_CH <sub>3</sub>	н	снз	Н	н
A732	СН3-	CH <sub>3</sub> H <sub>3</sub> C-√∑-{	н.	снз	Н	н
A733	СН3-	H <sup>2</sup> C CH <sup>2</sup>	Н	СНЗ	н	н
A734	CH3-	CH <sub>3</sub>	Н	СНЗ	н	н
A735	СН3-	H <sub>3</sub> C H <sub>3</sub> C	н	снз	Н	н
A736	СН3-	H <sub>3</sub> C	н	снз	н	н
A737	снз-	<b>F</b> - <b>F</b>	Н	снз	н	н
A738	снз-	F-{=}-}	н	СНЗ	н	н
A739	снз-	\$-	н	снз	Н	Н
A740	СН3-	€;	н	снз	н	Н
A741	CH3-	F	н	снз	Н	н
A742	CH3-		н	снз	Н	н
A743	СН3-	CI_CI	н	СНЗ	н	н
A744	СН3-	CI—⟨Ç]	н	CH3	н	Н
A745	снз-		Н	СНЗ	н	н
A746	снз-		Н	снз	н	н
A747	СН3-	CI—	н	СНЗ	н	н
A748	СН3-		н	снз	н	н
A749	СН3-	H <sub>3</sub> CO_OCH <sub>3</sub>	н	снз	Н	н
A750	снз-	H3CO-()-3	Н	снз	Н	н
A751	СН3-	H <sub>2</sub> CO	н	снз	н	н

No.	R1	R2	R3_	R4	R5	IR6
.,,,,,	<del>                                     </del>	OCH <sub>3</sub>		<u> </u>		
A752	CH3-	OCH;	Н	СНЗ	Н	н
A753	СН3-	300 /	н	СНЗ	Н	н
A754	СН3-	H <sub>3</sub> CO	н	СНЗ	н	н .
A755	СН3-	F_OCH <sub>3</sub>	Н	СН3	Н	н
A756	снз-	). Z '	н ·	снз	н .	н
A757	CH3-		Н	СНЗ	н	н
A758	снз-	OCH <sub>3</sub>	H .	СНЗ	н	н
A759	CH3-	©∺ F	H	снз	Н	Н
A760	СН3-	PCH <sub>3</sub>	Н	сн3	Н	н
A761	снз-	H <sub>3</sub> CO F—√□—i	н	снз	н	н
A762	снз-	H <sub>3</sub> CO	н	снз	н	н .
A763	снз-	H <sub>3</sub> CO_F	н	снз	Н	н
A764	снз-	H₃CO-{\(\bigcirc\)	н	снз	н	Н
A765	CH3-	H3CCC	н	снз	н	Н
A766	снз–	H <sub>3</sub> CO-	H·	снз	н	н
A767	снз-	CI_OCH <sub>3</sub>	н	СНЗ	Н	н
A768	СН3-	OCH₃ CI—⟨□}	н .	снз	Н	Н
A769	СН3-	©CH₃ CH3	н	СНЗ	Н	н
A770	снз-	ocH₃ G	н	снз	Н	н
A771	СН3	H₃CO CI— CI—	н	СНЗ	Н	н
A772	CH3-	H₃CO CI	н	СНЗ	н	Н
A773	CH3-	H₃CO_CI	н	СНЗ	Н	н

No.	RI	R2	R3	R4	R5	R6
A774	CH3-	H₃CO-⟨{}{}	Н	СН3	н	Н
A775	CH3-	H,CO CI	મુ	снз	н	н
A776	CH3-	CI, H₃CO-⟨□>—}	н	снз	Н	Н
A777	снз-	F_CH <sub>3</sub>	н	снз	н	н
A778	CH3-	CH <sub>3</sub> F—⟨∑}—;	н	СНЗ	н	Н
A779	CH3-	CH <sub>3</sub>	н	снз	н	Н
A780	снз-	CH <sub>3</sub>	н	снз	Н	н
A781	снз-	H <sub>3</sub> C F-\(\)	н	СНЗ	Н	н
A782	снз-	H <sub>3</sub> G F	Н	онз	н	н
A783	снз-	H <sub>3</sub> C_F	н	снз		н
A784	СН3-	H₃C-{	н	СНЗ	н	н
A785	CH3-	H <sub>3</sub> C +	н	СНЗ	Н	Н
A786	СН3-	H3C- <b>⟨</b>	н	СНЗ	н	н
A787	СН3-	Br. OCH <sub>3</sub>	н	снз	н	н
A788	CH3~	OCH₃ Br—{	н	СНЗ	н	н
A789	CH3-	Br DCH <sub>3</sub>	н	СНЗ	н	н
A790	СН3-	OCH₃ ⇒ Br	н	СНЗ	н	H
A791	снз-	H <sub>3</sub> CO Br	Н	СНЗ	н	н
A792	снз-	H <sub>3</sub> CO	Н	СНЗ	н	н
A793	снз-	H <sub>3</sub> CO_Br	н	снз	Н	н
A794	СН3-	H₃CO-{_}}-1	н	снз	н	н
A795	СН3-	H <sub>3</sub> CO Br	н	снз	Н	н

No.	Ri	R2	R3	R4	R5	R6
		Br				
A796	CH3-	H₃CO-(	H ·	CH3	Н	Н
A797	CH3-		н	СНЗ	н	н
A798	CH3-	OCH <sub>3</sub>	н	СНЗ	Н	Н
A799	снз-	ON-⟨S-OCH3	Н	СНЗ	н	н
A800	CH3-	H³CO }	н	снз	н	Н
A801	CH3-	H₃CO N-{_}	Н	СНЗ	н	н
A802	СН3-	Ch Ch	Н	СНЗ	Н	Н
·A803	снз-	-4	н	СНЗ	н	н
A804	снз-	POCHS F-C	н	СНЗ	н	н
A805	СН3-	H³C0-€ 1	н	СНЗ	Н	н
A806	СН3-	OCH <sub>3</sub> F—C≥→ OCH <sub>3</sub>	н	снз	н	н
A807	снз-	OCH3 OCH3	Н	СНЗ	н	Н
A808	снз-	CI CI	Н	снз	н	н
A809	снз-	CI-CS-1 CI	н	снз	н	н
A810	снз-	CI CI CI	н	снз	н	н
A811	СН3-	OCH <sub>3</sub>	н	снз	Н	н
A812	снз-	DCH <sup>2</sup> OCH <sup>3</sup>	н	снз	Н	н
A813	снз-	OCH3	н	СНЗ	н	н
A814	СН3-	H <sub>3</sub> CO	н	снз	Н	н
A815	снз-	H₃CO- <b>⟨_</b> }-{_}-{	н	СНЗ	н	н
A816	СН3-		Н	СНЗ	н	н
A817	СН3-	H <sub>3</sub> CO	н	СНЗ .	Н	Н

No.	R1	R2	R3	R4	R5	R6
A818	снз-	H300-(C)-(C)	н	снз	н	н
A819	снз-	OCH,	н	снз	н	н
A820	СН3-	H3C0	н	снз	н	н
A821	снз-	H₃CO-⟨>-{\}	н	снз	Н	н
A822	СН3-	<b>₫-</b>	н	снз	н	н
A823	СН3-	<b>₹</b> >-<>-,	н	снз	н	Н
A824	СН3-		н	снз	н	н
A825	онз-	<u>&amp;</u>	н	снз	н	н
A826	снз-		Н	СНЗ	н	Н
A827	СН3-		н	СНЗ	н	н
A828	снз-	Ø.	н	СНЗ	н	н
A829	снз-	<b>₽</b>	н	СНЗ	н	н
A830	CH3-		н	СНЗ	Н	н
A831	СН3-		н	СНЗ	Н	н
A832	снз-	CO'	н	СНЗ	Н	н
A833	СН3-	CT)-i	н	СНЗ	н .	н
A834	СН3-		н	СНЗ	н .	Н
A835	снз-		н	СНЗ	н	н
A836	СН3-	(Ch)	Н	СНЗ	н	н
A837	СН3-	,CO	н	снз	н	н
A838	СН3	Çî	н	СНЗ	н	н
A839	снз–	CC-1.	н	снз	н	н

No.	R1	R2	R3	R4	R5	R6
No.	100	74	173	17.4	פא	KO
A840	снз-		н	снз	н	н
A841	снз-	Č;	н .	снз	н	H
A842	СН3-	<b>'03</b>	н	СНЗ	н	н
A843	СН3-	,CC	н	СНЗ	н	н
A844	CH3~	Ţ?	н	СНЗ	н	Н
A845	CH3-	(I)	н	СНЗ	н	н
A846	CH3-		н	СН3	н	H
A847	CH3-		н	СНЗ	н	н
A848	CH3-	TOP	н	СНЗ	н	Н
A849	CH3-	,CC;	н	CH3	Н	н
A850	CH3-	ĈŢ\$	Н	снз	Н	н
A851	CH3-	CT <sub>p</sub>	Н	СНЗ	н	Н
A852	CH3-		н	снз	Н	Н
A853	CH3~	TOP	н	СНЗ	Н	H
A854	CH3~	,CT	н	СНЗ	н	н
A855	СН3-	Ţħ	н .	СНЗ	н	н
A856	СН3-	(CIN→	Н	СН3	н	Н
A857	СН3-	Ç <sup>1</sup>	н	СНЗ	Н	Н
A858	CH3-	TON N	Н	СНЗ	н	н
A859	CH3-		н	СНЗ	н	н
A86D	СН3-		н	снз	н	Ĥ

No.	R1	R2	R3_	R4	R5	R6
A861	CH3-	T,	н	снз	н	н
A862	СН3-	, CTN	Н	СНЗ	Н	н
A863	СН3-	<b>P</b> 3	н	СНЗ	Н	н
A864	СН3-	(Is)-+	Н	СН3	н	н
A865	СН3-	Ţ <sup>N</sup> S	н	СНЗ	н	н
A866	СН3-	TO S	н	снз	Н	н
A867	СН3-	, O's	н	снз	н	н
A868	снз-	ÇT°s	н	СНЗ	н	н
A869	снз-		Н	снз	Н	H
A870	СН3-		н	снз	н	н
A871	СН3-	TOW	н	снз	н	н
A872	CH3-	,CT%	н	снз	Н	н
A873	СН3-	Ĉ.	Н	снз	н	н
A874	снз-	OT, N	н	снз	н	н
A875	снз-		н	СНЗ	н	н
A876	CH3-	TOI;N	н	снз	Н	н
A877	СН3-	, LTgh	Н	снз	н	н
A878	CH3-	QT3N	н .	снз	н	Н
A879	СН3-	Ţ?	н	снз	н	н
A880	СН3-	,CC	н	снз	н	н
A881	СН3-	TOR	н	СНЗ	н	н
A882	СН3-	Č;	н	снз	н	н

No.	R1	R2	R3	R4	R5	R6
A883	СН3-	СН3-	н	Qu	н	н
A884	CH3-	СНЗСН2-	н	Qu	н	н
A885	снз-	<b>∕</b> ∕\	н	Qu	н	н
A886	снз-	Y	Н	Qu	н	Н
A887	СН3-	<b>√</b>	н	Qr	н	н
A888	СН3-	人、	н	Q	н	н
A889	СН3-	$\gamma$	н	Qi	н	н
A890	CH3-	7	н	Qu	Н	н
A891	снз-	~~\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	Qu	н	н
A892	снз-	Y~~	н	Qu	Н	н
A893	CH3-	X	н	Q	н	н
A894	СН3-	~	н	Qu	Н	Н
A895	CH3-	<b>\\\\</b>	н	Qu	н	н
A896	СН3-	<del></del>	н	Qu	Н	н
A897	СН3-	~~~`\	н		н	н
A898	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н		н	Н
A899	CH3-	n-C8H17-	Н		Н	н
A900	СН3-	人 <b>~~</b> ~	Н	Qu	н	н
A901	СН3-	Q	Н	Qu	Н	н
A902	СН3-		н	Qu	н	н
A903	CH3-		н	Qu	Н	Н
A904	снз-	$\triangleright \rightarrow$	н	Qu	Н	н

No.	R1	R2	R3	R4	R5	R6
A905	снз–	$\Diamond$ -1	н		Н	н
A906	CH3-	$\bigcirc\!$	н		Н	н
A907	CH3-	$\bigcirc$	н		Н	н
A908	CH3-	$\bigcirc$ -1	н		н	н
A909	CH3-	$\bigcirc$ -1	н		н	н
A910	CH3-		н		н	н
A911	онз-		н		н .	Н
A912	онз-	<b>-</b>	н		н	н
A913	СН3-	<u></u>	Н		н	н
A914	СН3-	F-(	н		н	н
A915	СН3-	F-(){	н		н	н
A916	снз–	F—C)m4	н		н	н
A917	СН3-	CI	Н		н	н
A918	СН3-	CI.	Н		н	н
A919	СН3-	c⊢ <b>(</b> )→(	н		н .	н
A920	СН3-	CH	н		н	н
A921	СН3-	C⊢ <b>(</b> )™	н		н	н
A922	снз-	Br	н		н	н
A923	СН3-	Br. —∤	Н		Н	н
A924	СН3-	Br <b>⟨</b> }-{	н.	Qu	Н	н
A925	CH3-	Br—()	н	Q	Н	н
A926	СН3-	Br—€∑i…{	н	Q	Н	Н

No.	R1	R2	lR3	R4	R5	R6
A927	снз-		н	Q	Н	H
A928	СН3-		Н	Q	н	Н
A929	СН3-		н	Q	н	Н
A930	СН3-	CH₃	н	Qu	н	Н
A931	СН3-	H <sub>3</sub> C	н	Qu	н	Н
A932	сӊз-	H₃C-{_}	н	Q	н	Н
A933	СН3-		Н	Q	н	н
A934	СН3-	n-C <sub>3</sub> H <sub>7</sub> -{{}}{	н		н	н .
A935	снз-	n-C <sub>4</sub> H <sub>9</sub> —{}	н	Q	н	н
A936	снз-	OH OH	Н		н	Н
A937	СН3-	HO	Н	Qu	н	н
A938	СН3-	HO-{}-{	н	Q	н	н
A939	СН3-	OCH₃	н	Q	Н	н
A940	СН3-	H <sub>3</sub> CQ	Н		Н	н
A941	CH3-	H3CO-	н		H .	Н
A942	СН3-	H3CO-	Н	Qu	Н	Н
A943	СН3-	H <sub>3</sub> CO-{_}\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	Q	н	н
A944	СН3-	OC <sub>2</sub> H <sub>5</sub>	Н	Q	н	Н
A945	СН3-	C <sub>2</sub> H <sub>5</sub> Q	Н	Q	н	н
A946	снз-	C <sub>2</sub> H <sub>5</sub> O-{}	н	Qu	н	Н
A947	снз-	n-C <sub>3</sub> H <sub>7</sub> O-{_}-{	Н		Н	н
A948	СН3-	n-C <sub>4</sub> H <sub>9</sub> O-{_}-{	Н	Q.,	н	н

No.	R1	R2_	R3	R4	IR5	R6
.,,,,		NO <sub>2</sub>			<del>                                     </del>	
A949	CH3-		Н		Н	Н
A950	CH3-	O <sub>2</sub> N	н		н	н
A951	снз-	O <sub>2</sub> N-{_}{	н	Q	н	Н
A952	СН3-	CN →	Н		н	н
A953	снз-	NC	н		н	н
A954	СН3-	NC-	Н		н	Н
A955	снз-	CF <sub>3</sub>	н		н	н
A956	СН3-	F <sub>3</sub> C △ →	н		н	н
A957	CH3-	F <sub>3</sub> C-{	н	Qu	н	Н
A958	снз-	COOH	н	Qu	н	н
A959	снз-	HOOC	н	Q	н	Н
A960	снз-	HOOC-{}	н		н	н
A961	СН3-	CO₂Me	Н		н	н
A962	снз-	MeO <sub>2</sub> C	н		н	н
A963	снз-	MeO <sub>2</sub> C-{}	н		Н	Н
A964	СН3-	CO <sub>2</sub> Et	н		н	H
A965	СН3-	EtO <sub>2</sub> C	Н		н	н
A966	снз-	EtO <sub>2</sub> C-\_\	н	Q	Н	н
A967	снз-	SMe ←	н		н	н
A968	снз-	MeS	н		н	н
A969	снз-	MeS-{	Н	Q	Н	н
A970	СН3-	SO₂Me	н	Q	н	н

No.	R1	R2	R3	R4	R5	JR6
		MeO <sub>2</sub> S	Ţ ·			- 1.0
A971	СН3-		Н		Н	н
A972	снз-	MeO <sub>2</sub> S-{}-{	Н	Q	Н	н
A973	снз-	NH₂ → ·	н		н	н
A974	СН3-	H <sub>2</sub> N	н		н	Н
A975	снз–	H <sub>2</sub> N-\_	н	Q	н	н
A976	снз-	NMe <sub>2</sub>	н	Q	н	н
A977	онз–	Me <sub>2</sub> N →	н	Q	н	н
A978	снз-	Me <sub>2</sub> N-(	н	Q	н	н
A979	снз-		н	Q	н	н
A980	СН3-		н	Q	н	н
A981	снз-		Н	Q	н	н
A982	снз-		н	Q	н	н
A983	снз-	Ov-Q	н	Q	н	н
A984	снз-	N-(2)-1	н	Q	н	н
A985	снз-	N-√	н	Q	н	н
A986	снз-		н	Q	н	н
A987	снз-	o_n-<>-	н	Q	н	н
A988	снз-	H³CN_N-⟨	н	Qu	н	н
<b>8989</b>	СН3-	H₃CN N-	н	Q	н	н .
A990	СН3-	H³CN_N-{}-!	н	Qu	н	н
A991	снз-	H <sub>3</sub> C_CH <sub>3</sub>	н	Q	н	Н
A992	снз-	сн <sub>3</sub> н₃с-⟨у;	н	Our	н	н

No.	R1	R2	R3	R4	R5	R6
A993	снз-	H³C CH²	н	Q	н	н
A994	снз-	CH <sub>3</sub> CH <sub>3</sub>	н	Q	н	н
A995	снз-	H <sub>3</sub> C H <sub>3</sub> C-⟨□}}	н	Qr	н	н
A996	СН3-	H³C .	н	Q	н	Н
A997	CH3-	<b>-</b>	н	Qu	н	н
A998	снз-	F-{\rightarrow}	н	Qu	н	н
A999	CH3-	<b>\$</b> -	н	Qu	н	н
A1000	СН3-	C.	н	Qr	н	н
A1001	снз-	F	Н		н	Н
A1002	снз-		н	Q	н .	н
A1003	снз-		н		Н	H
A1004	CH3-	CI CI	н		н	н
A1005	СН3-	G CI	н	Q	н	Н
A1006	СН3-	S.	н		н	н
A1007	снз-		Н		н	н
A1008	СН3-		н		н	Н
A1009	СН3-	H <sub>3</sub> CO_OCH <sub>3</sub>	н		н	н
A1010	СН3-	OCH <sub>3</sub> H <sub>3</sub> CO-⟨□⟩;	н	Q	Н	н
A1011	CH3-	OCH <sub>3</sub>	н	Qu	Н	н
A1012	СН3-	OCH <sub>3</sub>	н	Qu	н	н
A1013	CH3-	H₃CO H₃CO	н	Qu	н	Н
A1014	СН3-	H3C0	н	Q	н	н

No.	R1	R2	R3	R4	R5	R6
A1015	снз-	F_OCH <sub>3</sub>	н	Q	н	н
A1016	СН3-	OCH <sub>3</sub>	н	Qu	н	н
A1017	СН3-	OCH₃ F-€>	н	Qu	н	н
A1018	снз-	OCH <sub>3</sub>	н	Qu	н .	н
A1019	СН3-	OCH <sub>3</sub>	н	Qu	н	н
A1020	СН3-	OCH <sub>3</sub>	н	Q	Н	Н
A1021	СН3-	H <sub>3</sub> CO	н	Qu	н	н
A1022	СН3-	H3CO	н		н	н
A1023	снз-	H₃CO_F	н	Qu	н	н
A1024	снз-	H₃CO-{	н.	Q	н	н
A1025	снз-	H,CO H	н	Q	н	Н
A1026	снз-	H₃CO-⟨¬	н	Q	н	н
A1027	СН3-	CI_OCH <sub>3</sub>	н	Q	н	н
A1028	снз-	CI—(S)→	Н	Qi	н	Н
A1029	снз-	OCH <sub>3</sub>	н	Q	н	н
A1030	снз-	©CH <sub>3</sub>	н	Q.	н	н
A1031	снз-	CI—	н		н	н
A1032	снз-	H,CQ	н	Ou	н	Н
A1033	снз-	H₃CO_CI →	н		н	н
A1034	снз-	H₃CO-{\(\begin{array}{c} \text{CI} \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\ \\	н	Qu	н	н
A1035	снз-	H <sub>3</sub> CO	н		. Н	н
· A1036	СН3-	H³CO-{}	н	Q	н	н

No.	R1	R2	R3	R4	R5	R6
A1037	снз-	F_CH <sub>3</sub>	н	Or	н	н
A1038	сн3-	CH <sub>3</sub> F—⟨	Н	Q	н	н
A1039	CH3-	<b>₽</b>	н	Q	н	Н
A1040	CH3-	ÇH3	н		н	н
A1041	снз-	H <sub>3</sub> C F-\{	Н		Н	Н
A1042	снз-		н	Qu	н	н
A1043	СН3-	H₃C_F →	н	Q	н	н
A1044	снз-	H <sub>9</sub> C-⟨=⟩	н	Q	н	н
A1045	CH3-	4,c	н	Q	н	н
A1046	CH3-	H₃C-\	н	Q	н	н
A1047	СН3-	Br. OCH₃	н	Q	Н	н
A1048	CH3-	OCH <sub>3</sub>	н	Q	н	н
A1049	CH3-	Br OCH <sup>3</sup>	н	Q	н	Н
A1050	CH3-	OCH <sub>3</sub>	н	Q	н	н
A1051	СН3-	H₃CO Br—∰	н	· Ox	Н	н
A1052	СН3-	H,CO B	н	Ou	н	н
A1053	снз-	H <sub>3</sub> CO_Br	н		н	н
A1054	СН3-	H₃CO-⟨S)—I	н		н	н
A1055	CH3-	H <sub>3</sub> CO Br	Н	Q	н	н
A1056	снз-	H <sub>3</sub> CO-	н	Or	н	н
A1057	CH3-	H <sub>3</sub> CQ \	н	Q	Н	н
A1058	снз-	CN-€	н		н	н

No.	R1	R2	R3	R4	R5	R6
A1059	CH3-	N-{>-och3	н	Q	н	Н
A1060	снз-	H <sub>3</sub> CO >	н		Н	Н
A1061	СН3-	H <sub>3</sub> CO N	н	Q	н	н
A1062	СН3-	Ch.	н		н	н
A1063	снз-	F-C	н		н	Н
A1084	снз-	F-C-5 F	Н	Qu	н	н
A1065	СН3-	H3CO-(\$\frac{1}{2}\)	н	Qu	н	н
A1066	снз-	OCH <sub>3</sub>	н	Q	н	н
A1067	CH3-	осн³ осн³	н	Q	н	н
A1068	СН3-	ci—Ci	Н	Q	н	н
A1069	СН3-	CI—COCH3	н	Q	н	н
A1070	CH3-	H³CO-⟨∰t	Н		н	н
A1071	снз-	CH- CH3 CH3	н		н	н
A1072	снз-	н³со-⟨_}-} осн²	н		н	н
A1073	снз-	OCH <sub>3</sub>	н		н	Н
A1074	СН3-	H <sub>3</sub> CO	н		н	н
A1075	СН3-		н		н	н
A1076	СН3-	OCH <sub>3</sub> }:	н	Q	н	н
A1077	СН3-	H <sub>3</sub> CO ,	н		н	Н
A1078	CH3-	H3CO-{	н		н	Н
A1079	СН3-		н	Q	н	н
A1080	снз-		Н	Q	Н	н

No.	R1	R2	R3	R4	R5	R6
		H <sub>3</sub> CO-{_}-{_}				
A1081	CH3-	, , , , , , , , , , , , , , , , , , ,	H		H	Н
A1082	снз-	₫-O-1	н	Q	н .	Н
A1083	снз-	<u></u>	н	Q	Н	Н
A1084	СН3-	F-()-()-1	Н		Н	н
A1085	СН3-	₫-ď	Н		н	н
A1086	СН3-		н		Н	н
A1087	СН3-		н		н	Н
A1088	CH3-	4	н		н	н
A1089	СН3-	\$ <del>-</del> \$	Н	Qu	н	н
A1090	CH3-		н		н	н
A1091	CH3-		н		н	Н
A1092	СН3-	CCT'	н		Н	н
A1093	СН3-		н		Н	Н
A1094	СН3-		н	Q	н	н
A1095	СН3-		н	Q	Н	н
A1096	СН3-	TOT)	н	Q	н	н
A1097	СН3	,CT)	н	Q	н	н
A1098	онз-	Ç,	н	Q	н	н
A1099	онз-		н		н .	Н
A1100	CH3-	CT.	н		Н	Н
A1101	СН3-	Č	н	Q	н	н
A1102	СН3-	(CC)	н	Q	н	н

No.	R1	IR2	R3	R4	R5	R6
	1	6			11.0	110
A1103	CH3-	12 d	н		н	Н
A1104	снз-	Ç;	н	Qu	Н	н
A1105	снз~	CT\$-1	н	Qu	н	н
A1106	СН3-		Н	Qu	н	н
A1107	CH3-	Č:	н	Qr	н	н
A1108	CH3~	TOP	н	Qu	Н	н
A1109	CH3-	,CTS	н	Q	н	н
A1110	снз-	ÇT\$	Н	Q	Н	н
A1111	онз-		Н	Q	Н	Н
A1112	СН3-		н	Qi	н .	Н
A1113	CH3-		н	Qu	н	Н
A1114	CH3-	, CT'N	н	Qu	н	н
A1115	CH3-	ÇÜ	н	Qu	н	н
A1116	СН3-	CTN-i	н	Qu	н	н
A1117	CH3-	Č,	H	Q.	н	н
A1118	CH3-	, CIN	н	Q	н	Н
A1119	снз-	©l°+	н	Qu	н	н
A1120	CH3-	Č';	Н	Qi	Н	н
A1121	CH3-	'CIN	н	Qu	Н	н
A1122	СН3-	, CIN	н	Q	H	н
A1123	CH3-	Ţ.)	Н	Q	н .	н
A1124	СН3-	I S	н	Q	Н	н

No.	R1	R2	R3	R4	R5	R6
A1125	CH3-	, N	н	Q	H	н
A1126	снз-	TIN	н	Q	н	Н
A1127	снз-	, CI,	н	Q	н	н
A1128	снз-	T <sub>s</sub>	н	Q	н	н
A1129	снз-		н		Н	н
A1130	СН3-	Ĉ()	н		н	н
A1131	СН3-	"CO"	н		Н	н
A1132	CH3-	, OG	н		н	н
A1133	CH3-	Ţ,	н		н	н
A1134	СН3-		н		н	н
A1135	СН3-	Ţ,	Н		н	Н
A1136	онз-	TO W	Н	Qu	н	н
A1137	СН3-	,CT3N	н		Н	н
A1138	CH3-	Ĉ.	н	Qu	н	н
A1139	CH3-	Ç.	н	Q	н	н
A1140	CH3-	,CC	н		н	н
A1141	СН3-	TCT)	н	Qu	н	н
A1142	CH3-	Ĉ,	н		н	н
A1143	снз-	СН3-	н	l,	н	н
A1144	CH3-	GH3CH2-	н	گ,	н	Н
A1145	СН3-	<b>^</b> ``	н	l,	н	Н
A1146	снз-	Y	н	2,	Н	н

No.	R1	R2	R3	R4	R5	R6
A1147	снз-	<b>\\\</b>	н	Å,	н	н
A1148	снз-	人、	н	Ŷ,	н	н
A1149	GH3-		н	Ŷ,	н	н
A1150	СН3-	丫	Н	l,	н	н
A1151	снз-	~~``\	н	گ <sub>ه</sub>	н	н
A1152	снз-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	<u></u>	н	н
A1153	СН3-	<u> </u>	н	Å,	н	н
A1154	снз-	7	Н	<u>گ</u> ہ	Н	н
A1155	СН3-	~~~	н	l,	Н	н
A1156	снз-	<del>                                      </del>	н	ئى ر	Н	н
A1157	СН3-	~~~``\	н	l,	н	н
A1158	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	<u>گ</u>	н	н
A1159	СН3-	n-C8H17-	н	Ŷ,	н	н
A1160	СН3-	L~~~	н	گہ	н	н
A1161	CH3-	Qu	н	l,	н	н .
A1162	CH3-		н	l,	н	Н
A1163	CH3-		Н	١,	н	н
A1164	снз-	<b>&gt;</b> →	н	<u>گ</u>	н	н
A1165	снз-	$\Diamond$ -1	н	گ,	н	н
A1166	снз-	$\bigcirc$	н	l,	н	н
A1167	снз-	$\bigcirc$ -1	н	گ <u>ہ</u> گہ	н	н
A1168	снз-		н	الم الم	н	н

No.	R1	R2	R3	R4	R5	R6
A1169	снз-	<b>◯</b> -₁	н	<u> </u>	н	н
A1170	CH3-		н	<u>گ</u>	н	н
A1171	CH3-	<b>⊘</b> m4	Н	l,	н	Н
A1172	СН3 <del></del>	F	н	l,	н	Н
A1173	CH3-	F	н	<u>گ</u> ر	н	н
A1174	снз-	F-(	н	Ŷ,	н	H
A1175	снз-	F-(>-1	н	<u>گ</u>	н	н
A1176	снз-	F———	н	١,	н	н
A1177	СН3-	CI →	н	2,	н	н
A1178	СН3-	CI.	н	Î,	Н	н
A1179	CH3-	c⊢ <u></u>	н	Ŷ,	Н	н
A1180	снз-	ci—(i	н	ي ا	Н	н
A1181	CH3-	CI—(	н	<u>گ</u>	н	н
A1182	CH3-	Br	н	٩	н	н
A1183	СН3-	Br.	н	Ŷ,	н	н
A1184	снз-	Br—⟨{	н	ئى ك	н	н
A1185	СН3-	Br—{	Н	Ŷ,	н	н
A1186	СНЗ-	Br─€∑™{	н	l,	н	н
A1187	CH3-		н	گي	Н	н
A1188	СН3-		н	2,	Н	Н
A1189	СН3-	<del> </del>	н	l,	Н	н
A1190	снз–	CH <sub>3</sub>	н	1,	н	н

No.	R1	R2	R3	R4	R5	R6
NO.	121	H <sub>3</sub> C	I/O	0	INO .	1,70
A1191	СН3-	· · ·	н	ير .	н	H
A1192	СН3-	H <sub>3</sub> C-{	Н	Å,	Н	н
A1193	СН3-		Н	l,	н	н
A1194	СН3-	n-C <sub>3</sub> H <sub>7</sub> -	н	l,	н	н .
A1195	CH3-	n-C <sub>4</sub> H <sub>9</sub> {}-{	н	ئى ا	н	н
A1196	CH3-	OH →	Н	ئى ا	н	н
A1197	CH3-	HO ———	н	یا <sub>ن</sub>	Н	Н
A1198	CH3-	HO-{_}-{	Н	گ <sub>ا</sub>	Н	н
A1199	CH3-	OCH₃	Н	Ŷ,	Н	н
A1200	CH3-	H₃CO <u></u> —∤	Н	برگ	H <sub>.</sub>	н
A1201	снз-	H₃CO- <b>⟨</b> }-	н	بُ	н	Н
A1202	снз-	H <sub>3</sub> CO-	н	Ŷ,	н	н
A1203	снз-	H <sub>3</sub> CO-{\bigs_m\}	н	٩	H	н
A1204	снз-	OC <sub>2</sub> H <sub>5</sub>	Н	١	н	н
A1205	снз-	C <sub>2</sub> H <sub>5</sub> Q	н	Ŷ,	н	н
A1206	онз-		н	ئى ر	Н	Н
A1207	СН3-	n-C <sub>3</sub> H <sub>7</sub> O-{_}-{	н	ئہ	н	н
A1208	снз-	n-C <sub>4</sub> H <sub>9</sub> O-{_}-{	н	l,	н	н
A1209	снз-	NO <sub>2</sub>	H	ئى _	н .	н
A1210	снз-	O <sub>2</sub> N	Н	ئى ر	н	н
A1211	СН3-	02N-(	н	بُ	н	н
A1212	СН3-	CN C	н	ئى,	н	н

No.	R1	R2	R3	R4	R5	R6
A1213	СН3-	NC	н	الم الم	н	н
A1214	снз-	NC-{_}-{	н	Ů,	н	Н
A1215	СН3-	CF <sub>3</sub>	н	١	н	Н
A1216	СН3-	F <sub>3</sub> C	н	2,	н	н
A1217	СН3-	F <sub>3</sub> C-{}-{	н	l,	Н	н
A1218	СН3-	соон	н	بُ	н	н
A1219	СН3-	HOOC	н	ئى ر	н	н
A1220	снз-	H00C-{_}-	н	گہ	н	н
A1221	СН3-	CO₂Me	н	Ŷ,	н	н
A1222	СН3-	MeO <sub>2</sub> C	н	Î,	н	н
A1223	снз-	MeO <sub>2</sub> C-{}	н	Î,	н	н
A1224	СН3-	CO <sub>2</sub> Et	н	<u></u>	н	Н
A1225	СН3-	EtO <sub>2</sub> C	н	ئ,	н	н
A1226	CH3-	EtO <sub>2</sub> C-{}	н	Ŷ,	н	Н
A1227	снз-	SMe	н	گہ	н	н
A1228	снз-	MeS	н	Ů,	Н	н
A1229	СН3-	MeS-{_}	н	<u>گ</u>	Н	н
A1230	СН3-	SO₂Me	н	ئے	Н	Н
A1231	снз-	MeO <sub>2</sub> S	н	ئى ر	н .	Н
A1232	снз-	MeO <sub>2</sub> S-{}	н	l,	Н	н
A1233	СН3-	NH₂ →	н	Î,	н	н
A1234	снз-	H <sub>2</sub> N	н	گې	н	н

No.	R1	IR2	R3	R4	R5	R6
A1235	снз-	H <sub>2</sub> N-\\\\\\\	Н	Ŷ,	н	н
A1236	снз-	NMe <sub>2</sub>	н	2,	н	н
A1237	CH3-	Me <sub>2</sub> N	н	بُ	н	н
A1238	снз-	Me₂N-⟨¯}—ţ	н	Ŷ,	н	н
A1239	CH3-	\(\rac{1}{2}\rightarrow\)	н	گي.	Н	Н
A1240	снз-		н	بُر	н	н
A1241	снз-	Cn-<>-;	Н	گ <sub>ا</sub>	н	Н
A1242	снз-		н	Ŷ,	н	н
A1243	снз-	O-Q	н	Î,	н	н
A1244	снз–	O₁-(C)-1	н	l,	Н	Н
A1245	СН3-	O_N-{\rightarrow}	н	بگ	Н	Н
A1246	CH3-	ON-⟨∑	н	l,	Н	н
A1247	CH3-	O_N-< <u></u> _}-	н	l,	Н	н
A1248	СН3-	H3CN N-	н	l,	Н	н
A1249	CH3-	H3CN N-	н	l,	н	н
A1250	сн3-	H3CN N-	н	ا ا	н	н
A1251	снз-	H <sub>3</sub> C_CH <sub>3</sub>	н	l,	Н	н
A1252	CH3	H₃C-⟨∑){	Н	l,	Н	Н
A1253	снз-	H <sub>3</sub> C-M <sub>3</sub> CH <sub>3</sub> H <sub>3</sub> C	н ,	l,	Н	н
A1254	СН3-	CH <sub>3</sub>	н	上, 上,	н	н
A1255	СН3-	H₃C H₃C-⟨¯}—ţ	н	l,	н	н
·A1256	CH3-	H <sub>3</sub> C -1	н	l,	н	н.

No.	R1	R2	R3	R4	R5	R6
A1257	СН3-	F_F 	н	Ŷ,	н	H .
A1258	снз-	F-{\(\sigma\)_+}	н	l,	Н	н
A1259	СН3-	Ç,	н	<u>گ</u>	Н	н
A1260	СН3-	¢	н	<u>گ</u>	Н	Н
A1261	CH3-	F—————————————————————————————————————	н	<u>گ</u> ر	Н	н
A1262	CH3-	<b>\</b>	н	<u>گ</u> ر	н	н
A1263	CH3-		н	Ŷ,	н	н
A1264	СН3-	CI—(□)—;	н	Î,	Н	Н
A1265	СН3-		н	Ŷ,	н	Н
A1266	СН3-	or Cartesian (Cartesian Cartesian C	н	Ŷ,	н	н
A1267	СН3-		н	Ŷ,	Н	н
A1268	СН3-		н	<u>گ</u>	н	Н
A1269	СН3-	H <sub>3</sub> CO OCH <sub>3</sub>	н	ئى ر	н	н
A1270	снз-	H <sub>3</sub> CO-	н	ئى ا	н	н
A1271	снз-	H²CO ⟨□⟩→ OCH²	н	کہ	Н	н
A1272	снз-	OCH3	н	Ŷ,	н	н
A1273	CH3-	H₃CO_ H₃CO-	н	<u>گ</u> ر	Н	н
A1274	СН3-	H <sub>2</sub> CO	н	١,	н	н
A1275	СН3-	F_OCH <sub>3</sub>	н	١,	н	н
A1276	CH3-	OCH <sub>3</sub>	н	يا,	Н	н
A1277	CH3-	OCH <sub>3</sub>	н	بُ	н	Н
A1278	CH3-	OCH <sub>3</sub>	н	Ů,	н	н

No.	R1	R2	R3	R4	R5	R6
A1279	онз-	J <b>e</b> f ∣	н	Ŷ,	н	н
A1280	СН3-	F	н	l,	н	Н
A1281	CH3-	H₃CO F—(_)—;	н	l,	Н	н
A1282	снз-	H₃CQ F	н	l,	н	н
A1283	снз-	H <sub>3</sub> CO_F	н	l,	н	н
A1284	снз-	H₃CO-{	Н	<u></u>	н	н
A1285	снз-	H³C0 H³C0	н	گ <sub>ا</sub>	н	н
A1286	снз-	н₃со-С	Н	گ <sub>ا</sub>	Н	н
A1287	снз-	CI_OCH <sub>3</sub>	н	l,	н	н
A1288	снз-	OCH <sub>3</sub>	н	Î,	Н	н
A1289	снз-	ocH₃ G	Н	Å,	н	н
A1290	снз-	CI CI	н	گ <sub>ن</sub>	н	н
A1291	снз-	CI—(	н	٨	н	н
A1292	снз-	CI → H³CO	н.	Ŷ,	Н	н
A1293	CH3-	H <sub>3</sub> CO_CI	н	l,	Н	Н
A1294	СН3-	H³CO-{	Н	Ŷ,	н	н
A1295	СН3-	H³cco Ci	н	l,	н	н
A1296	СН3-	H³CO-⟨	н	<u></u>	Н	н
A1297	СН3-	F_CH <sub>3</sub>	н		н	н
A1298	снз-	CH <sub>3</sub> F—{}	н	گې	н	н
A1299	снз-	CH₃ F	н	<u></u> <u>L</u> ,	н	н
A1300	снз-	F CH3	н	گ,	н	н

No.	R1	R2	R3	R4	R5	R6
A1301	CH3-	H <sub>3</sub> C F-(-){	н	<u></u>	н	н
A1302	CH3-	H <sup>2</sup> C +	Н	گ,	Н	н
A1303	снз-	H <sub>3</sub> C_F	н	بُر	н	н
A1304	снз-	H₃C-{\bigc}	н	<u>گ</u>	Н	Н
A1305	снз-	H <sub>3</sub> C	н	<u>ڳ</u>	Н	н
A1306	снз-	H <sub>3</sub> C-\	н	١	н	н
A1307	СН3-	Br. OCH₃	н	٩	н	Н
A1308	CH3-	осн₃ вг—б—∤	Н	Ů,	н	н
A1309	снз-	OCH <sub>3</sub>	Н	Ŷ,	н	н
A1310	СН3-	DCH <sub>3</sub> Br	н	گہ	Н	н
A1311	CH3-	H₃CO Br———}	H <sub>.</sub>	بُ	н	н
A1312	CH3	H <sub>2</sub> CO	н	<u></u>	н	Н
A1313	снз-	H₃CO_Br	H	کہ	н	н
A1314	снз-	H₃CO-⟨□⟩	н	1,	н	Н
A1315	СН3-	H <sub>3</sub> CO Br	н	Ŷ,	н	н
A1316	снз-	Br. H₃CO-⟨¯_}-	н	Î,	Н	н
A1317	СН3-	H <sub>3</sub> CO_>	н	١	н	н
A1318	снз-	OCH <sub>3</sub>	н	<u></u>	н	н
A1319	снз-	CN-⟨\$-ocH³	н	2,	Н	н
A1320	СН3-	H <sub>3</sub> CO >	н	2,	н	н
A1321	СН3-	H₃CO N-√	н	<u></u> <u>,</u>	н	н
A1322	CH3-	Ch O	н	l,	н	н

No.	R1	R2	R3	R4	R5	R6
A1323	снз-	F-\$	Н	ا ا	н	н
A1324	СН3-	F-C	н	<u>گ</u> ,	н	н
A1325	СН3	н³со-С\$-1	н .	ئہ	н	н
A1326	СН3-	OCH <sub>3</sub>	Н	ب	н .	н
A1327	CH3-	H³CO-⟨∑H DCH³	Н	l,	н	н
A1328	СН3-	a-(2)	н	l,	н	н
A1329	СН3-	CH-⟨□>+ CI	Н	ئى,	Н	н
A1330	СН3-	H3CO-(2)	н	ئى,	Н	Н
A1331	CH3-	CI-€>-  CI-€>-	н	<u>گ</u>	н	Н
A1332	CH3-	H³CO-{\rightarrow} H\$CO-{\rightarrow} H\$CO-{\righta	н	l,	н	Н
A1333	СН3-	OCH <sub>3</sub>	Н	<u></u>	н	н
A1334	СН3-	H <sub>3</sub> CO	н	Î,	н	Н
A1335	СН3-	H3CO-{}-{}-{	н	Î,	н	н
A1336	СН3-	OCH <sub>3</sub> };	н	٩	Н	н
A1337	снз-	H <sub>3</sub> CO	н	ئہ	Н	н
A1338	снз-	H <sub>3</sub> CO-()-() <sup>1</sup>	н	١	н	Н
A1339	СН3-	OCH.	н	٤	н	н
A1340	СН3-	H-3C0	н	ئى ا	Н	н
A1341	СН3-	H₃CO-(\$\)—	Н	٤	н	н
A1342	СН3-	Ø-0-	н	١	н	н
A1343	CH3-	<u></u>	н	١,	Н	н
A1344	CH3-	F-(-)-(-)-1	н	l,	н	н .

No.	R1	R2	R3	R4	R5	R6
A1345	СН3-	d-d`	Н	Î,	н	н
A1346	снз-		Н	l,	н .	Н
A1347	СН3-		н	l,	Н	н
A1348	CH3-	Q-\(\rho\)	Н	l,	н	Н
A1349	CH3-	\$\ldot\\ \O \\ \O	н	<u>گ</u>	Н	н
A1350	СН3-		н	Î,	Н	н
A1351	CH3-		н	گ <sub>ا</sub>	Н	н
A1352	CH3-	CO,	н	l,	н .	н
A1353	снз-		Н	i,	н	н
A1354	снз-		Н	<u></u>	Н	н
A1355	CH3-		н	<u>گ</u>	н	Н
A1356	CH3-	TON	н	Î,	н	Н
A1357	снз-	, DY	н	l,	н	Н
A1358	СН3-	<b>P</b>	Н	l,	н	Н
A1359	СН3-		н	l,	н	н
A1360	CH3-		н	l,	н	н
A1361	снз-	j.	н	گې	н	н
A1362	снз-	(CI)	н	l,	н	Н
A1363	СН3-	,CC	н	<u>L</u> ,	н .	н
A1364	CH3-	Ç.	н	<u></u>	н	н
A1365	снз-	(Is)	н	<u></u>	н	Н
A1366	CH3-		н	l,	Н	н

No.	R1	R2	R3	R4	R5	R6
110.	<u> </u>	17			113	110
A1367	СН3-	(C)	н	l,	н	Н
A1368	снз-	TOP	н	l,	н	н
A1369	снз-	,CI?	н	l,	Н	н
A1370	снз-	Ğ.	Н	L,	н .	н
A1371	снз-		Н	٩	н	н
A1372	CH3-		Н	Î,	н	н
A1373	СН3-	TON	н	Ŷ,	н	н
A1374	СН3-	, LTM	н	Ŷ,	н	н
A1375	СН3-	ĞÜ,	н.	Ŷ,	н	н
A1376	CH3-		н	2,	Н	Н
A1377	СН3-	Ç, Ç,	н	٤	н	н
A1378	CH3-	TO N	н	ب	Н	н
A1379	CH3-	(T)-1	н	<u>گ</u> ہ	Н	н
A1380	CH3-	Ţ,	н	گ, گ,	н	н
A1381	СН3-	'CI'	н	l,	н .	н
A1382	снз-	, CT)	н	Î,	Н	н
A1383	снз-	Ť.	н	ا ا	ļ	н
A1384	снз-	OL»-+	н	l,	н	н
A1385	снз-	ŌN S	н	l,	н	н
A1386	СН3-	TO NS	н	ئے	н	н
A1387	CH3-		н	<u></u> <u></u> ,	н	н
A1388	CH3-	T's	н	١,	н	Н

No.	R1	R2 ·	R3	R4	[R5	R6
A1389	СН3	Ci,	н	Ŷ,	Н	н
A1390	снз-	Č.	н	بُ	н	н
A1391	СН3-	TOW	н	l,	н	н
A1392	снз-	,CTD	Н	Ŷ,	н	н
A1393	СН3-	Ĉ.	Н	Ŷ,	н	н
A1394	CH3-		Н	Ŷ,	н	н
A1395	СН3	ĈŢ'n	н	ئى	н	н
A1396	СН3-	(CIE)	н	l,	н	н
A1397	СН3-	,CT3N	н	l,	н	н
A1398	снз-	ĈŢŝ'n	н	l,	Н	н
A1399	CH3-	Q:	н	کہ	н	н
A1400	CH3-	,CC	н	<u>گ</u>	н	Н
A1401	снз-	TOO;	Н	Ŷ,	н .	н
A1402	CH3-	Ğ;	Н	Ŷ,	Н	Н
A1403	снз-	снз-	СН3-	н	Н	н
A1404	снз-	CH3CH2-	CH3-	н	Н	н
A1405	CH3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	CH3-	н	н	н
A1406	СН3-	丫	снз-	н	н	Н
A1407	снз-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	снз-	н	н	н
A1408	снз-	L,	снз-	Н	н	н
A1409	CH3-	~~	снз-	Н	н	н
A1410	CH3-	丫	СН3-	Н	н	н

No.	R1	R2	R3	R4	R5	R6
A1411	снз-	$\wedge \wedge \lambda$	CH3~	Н	н	н
A1412	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	CH3-	н.	н	Н
A1413	CH3-	大ス	Сн3-	н	Н	н
A1414	снз-	7	СН3-	Н	Н	Н
A1415	СН3-	<b>\\\\\</b>	СН3-	н	н	Н
A1416	СН3-	人、	СН3-	Н	н	Н
A1417	снз-	<b>^</b>	СН3-	н	H	Н
A1418	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	СН3-	н	Н	н
A1419	CH3-	n-C8H17-	СН3-	Н	н	н
A1420	снз–	L~~~	CH3-	н	Н	н
A1421	СН3-	Qu	СН3-	н	H	н
A1422	CH3-		Сн3-	н	н	н
A1423	СН3-	Q	снз-	н	н	н
A1424	снз-	ightharpoonup	снз-	н	н	н
A1425	СН3-	$\Diamond$	снз-	н	н	н
A1426	снз-		СН3-	н	н	н
A1427	CH3-		СН3~	н	н	н
A1428	снз-	$\bigcirc \dashv$	СН3-	Н	н	н
A1429	СН3-	<u></u>	CH3-	н	н	н
A1430	СН3-		CH3-	н	н	н
A1431	снз-	<b>⊘</b> {	CH3-	н	Н	н
A1432	СН3-	F	СН3-	н	н	н

No.	R1	IR2	R3	R4	R5	R6
A1433	снз-	F				н
A1434	снз-	F-{}-{	снз–	н	н	н .
A1435	СН3	F-{_}-	СН3-	Н	Н	Н
A1436	СН3-	F-{>{	СН3-	Н	Н	Н
A1437	снз-	CI	СН3-	н	н	Н
A1438	снз-	CI	снз-	н	Н	н
A1439	снз-	CH	снз–	н	н	н
A1440	снз-	CI-()-1	СН3-	Н	Н	н
A1441	CH3-	CH	СН3-	н	Н	н
A1442	снз-	Br	СН3-	н	н	н
A1443	CH3-	Br.	СН3-	Н	Н	н
A1444	СН3-	Br—	СН3-	н	Н	н
A1445	СН3	Br—()	CH3-	Н	Н	Н
A1446	СН3-	Br———	СН3-	н	Н	Н
A1447	СН3-	<b>△</b>	CH3-	н	н	Н
A1448	СН3-	<u></u>	СН3-	н	Н	Н
A1449	снз-	<del>  _</del>	СН3-	н	н	н
A1450	CH3	CH₃ <	снз-	н	н	н
A1451	СН3-	H <sub>3</sub> C	снз-	н	н	н
A1452	СН3-	H <sub>3</sub> C-{{}}	снз-	н	н	н
A1453	СН3-		СН3-	н	н	н
A1454	снз-	n-C <sub>3</sub> H <sub>7</sub> —{}	снз-	н	н	н

No.	R1	R2	R3	R4	R5	R6
A1455	CH3-	n-C <sub>4</sub> H <sub>9</sub> —{	CH3-	н	н	Н
A1456	CH3-	ОН	CH3-	н	Н	н
A1457	CH3-	HO	CH3-	н	Н	н
A1458	СН3-	HO-{\bar{\bar{\bar{\bar{\bar{\bar{\bar	снз-	Н	Н	Н
A1459	снз-	OCH <sub>3</sub>	СН3-	Н	н	н
A1460	СН3-	H <sub>3</sub> CO	CH3-	Н	н	н
A1461	CH3-	H3CO-{}-{	CH3-	Н	н	н
A1462	снз-	H <sub>3</sub> CO-{}	снз-	н	н	Н
A1463	снз-	H <sub>3</sub> CO-{	СН3-	н	н	Н
A1464	СН3-	OC₂H <sub>5</sub>	CH3-	н	Н	н
A1465	снз-		СН3-	н	н	н
A1466	CH3-	C <sub>2</sub> H <sub>5</sub> O-{}	СН3-	н	н	н
A1467	CH3-	n-C <sub>3</sub> H <sub>7</sub> O-	CH3-	н	н	н
A1468	снз-	n-C <sub>4</sub> H <sub>9</sub> O-	CH3-	н	н	Н
A1469	СН3	NO <sub>2</sub>	CH3-	н	н	н
A1470	CH3-	O <sub>2</sub> N	CH3-	н	н	н
A1471	снз-	O <sub>2</sub> N-{}	снз-	н .	н	н
A1472	снз-	CN	снз-	н	н	н
A1473	снз-	NC.	снз-	Н	н	н
A1474	снз-	NC-{}	снз-	Н	н	н
A1475	СН3-	NH <sub>2</sub>	СН3-	н	Н	н
A1476	CH3-	H <sub>2</sub> N	СН3∹	н	н	н

No.	R1	R2	R3	R4	R5	R6
		H <sub>2</sub> N-\\\\\\\				1
A1477	СН3-		CH3-	Н	Н	Н
A1478	CH3-	NMe <sub>2</sub>	снз-	Н	н	н
		Me N				
A1479	снз-	Me <sub>2</sub> N →	CH3-	н	Н	н
A1480	СН3-	Me <sub>2</sub> N-(T)	CH3-	н	н	н
A1481	CH3-		СН3-	Н	н	Н
A1482	CH3-	(N-()	снз-	Н	н	н
A1483	CH3-	()+( <u>)</u> -1	CH3-	Н	н	н
A1484	CH3-		CH3-	Н	H.	Н
A1485	CH3-	Ov-Q_	СН3-	н	Н	Н
A1486	снз-		СН3-	н	н	Н
A1487	СН3-	<b>○</b> \-\_	снз-	Н	н	н
A1488	CH3-		СН3-	н	н	Н
A1489	СН3-	<b>○</b> \v <del>-</del> (>-i	СН3~	н	н	н
A1490	СН3-	H3CN_N-	онз-	н	н	н
A1491	СН3-	H3CN N-	снз-	н	Н	н
A1492	СН3-	H3CN N-()-1	снз-	н	<b>н</b>	н
A1493	СН3-	OCH <sub>3</sub> F—⟨□}—∤	CH3-	Н	Н	Н
A1494	СН3-	OCH <sub>3</sub>	СН3-	Н	н	н
A1495	СН3-	OCH <sub>3</sub> F—(	Сн3-	н	н	Н
A1496	СН3-		СН3-	н	н	н
A1497	снз-	CCC'	снз-	н	н	н
A1498	снз-	СН3-	H	н	СН3	н

No.	R1	R2	R3	R4	R5	R6
A1499			Н	Н	СН3-	н
A1455	Ons-					<u> </u>
A1500	снз-	<b>~</b> ~``	Н	н	CH3-	н
A1501	снз-	$\stackrel{\sim}{-}$	н	Н	онз-	Н
A1502	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	СН3-	Н
A1503	снз-	人、	н	Н	СН3-	Н
A1504	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	снз-	н
A1505	снз-	丫	н	н	снз-	н
A1506	СН3-	<b>~~</b> ~	н	н	снз-	Н
A1507	СН3-	<b>\_</b> `	н	н	снз-	н
A1508	СН3-	Xx	н	н	снз-	н
A1509	СН3-	7	н	н	снз–	н
A1510	снз-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	снз-	н
A1511	снз-	人~	н	н	снз-	н
A1512	снз-	^^^\	н	н	снз-	н
A1513	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	снз-	н
A1514	онз-	n-C8H17-	Н	Н	онз-	н
A1515	CH3-	L	н	Н	СН3-	н
A1516	СН3-	Qu	н	Н	снз-	Н
A1517	CH3-		н	Н	СН3-	н
A1518	CH3-	Q	н	H	СН3-	н
A1519	CH3-	ightharpoonup	н	н	снз-	н
A1520	CH3-	$\Diamond \dashv$	н	Н	сн3-	н

No.	R1	Ř2	R3	R4	R5	R6
A1521	снз-		н	н	СН3-	н .
A1522	снз-	$\bigcirc$ $\dashv$	Н	Н	CH3-	н
A1523	СН3-	$\bigcirc$ $\dashv$	н	Н	СН3-	Н
A1524	снз-	<b>△</b> -₁	Н	Н	СН3	н
A1525	снз-		Н	н	СН3-	Н
A1526	СН3-	<b>⊘</b> 1114	Н	н	СН3	н
A1527	CH3-	<b>5</b> 4	н	н .	онз-	н
A1528	CH3-	F	Н	Н	СН3-	Н
A1529	СН3-	F-(){	Н	H	СН3-	н
A1530	CH3-	F-(>-1	н	н	снз-	н
A1531	снз-	F——>iii4	н	н	СН3	н
A1532	CH3	CI →	н	Н	СН3-	н
A1533	СН3-	CI	н	Н	СН3-	н .
A1534	снз-	c <del>-</del>	н	Н	CH3-	Н
A1535	снз-	CI—()	Н	н	CH3-	H
A1536	снз-	CH	н	н	СН3-	Н
A1537	СН3-	Br	н	н	CH3-	н
A1538	снз-	Br.	н	н	CH3-	н
A1539	СН3-	Br-{}-{	Н	Н	СН3-	н
A1540	снз-	Br—{}	н	Н	снз-	н
A1541	СН3-	Br—{	н	н	снз-	Н
A1542	СН3-	<b>⟨</b> }-₁	н	н	снз-	н .

No.	R1	R2	R3	R4	R5	R6
A1543	CH3-	1,	н	Н	СН3-	н
A1544	CH3-		Н	н	снз-	Н
A1545	CH3-	CH₃	н	Н	снз-	Н
A1546	CH3-	H <sub>3</sub> C	Н	Н	снз-	н
A1547	снз-	H <sub>3</sub> C-\	н	н	СН3-	н
A1548	СН3-	C <sub>2</sub> H <sub>5</sub> —{}	Н	н	СН3-	Н
A1549	снз-	n-C <sub>3</sub> H <sub>7</sub> -	н	Н	СН3-	н
A1550	СН3-	n-C₄H <sub>9</sub> -{}	н	н	CH3-	н
A1551	СН3-	OH OH	н	н	СН3-	н
A1552	СН3-	HO	н	н	CH3-	н
A1553	CH3-	но-{}-	н	н	СН3-	н
A1554	СН3-	OCH₃	н	н	снз-	н
A1555	СН3	H₃CQ —}-{	н	н	СН3-	Н
A1556	СН3-	H₃CO-{_}}-{	н	н	СН3- '	н
A1567	CH3-	H <sub>3</sub> CO-{_>-1	н	н	СН3-	н
A1558	снз-	H₃CO-⟨⟩ı•∮	Н	н	СН3-	н
A1559	снз-	OC <sub>2</sub> H <sub>5</sub>	н	Н	онз-	н
A1560	CH3-	C <sub>2</sub> H <sub>5</sub> Q	н	н	СН3-	н
A1561	CH3-	C₂H₅O-{\rightarrow}-{\rightar	н	н	снз-	н
A1562	СН3-	n-C <sub>3</sub> H <sub>7</sub> O-	Н	н	СН3-	н
A1563	снз-	n-C <sub>4</sub> H <sub>9</sub> O-{{1}	н	н	снз-	н
A1564	онз-	NO <sub>2</sub>	н	н	снз-	н

No.	R1	R2	R3	R4	R5	R6
		R2 O <sub>2</sub> N				
A1565	CH3		Н	Н	СН3-	Н
A1566	снз-	O <sub>2</sub> N-{}	н	н	СН3-	н .
A1567	снз–	CN →	н	н	снз-	н
A1568	СН3-	NC	н	Н	СН3-	н
A1569	снз-	NC-{_}-I	н	н	снз-	н
A1570	снз	NH <sub>2</sub>	н	н	СН3-	н
A1571	СН3-	H <sub>2</sub> N	Н	н	снз-	н
A1572	снз-	H <sub>2</sub> N-{}	н	н	снз-	н
A1573	CH3-	NMe <sub>2</sub>	н	Н	снз-	H
A1574	СН3-	Me <sub>2</sub> N	н	Н	снз-	н
A1575	снз-	Me <sub>2</sub> N-⟨¯⟩;	н	Н	СН3-	н
A1576	СН3-		н	Н	СН3-	Н
A1577	СН3-	Cn-€}	н	Н	снз-	н
A1578	снз-	Cn-⟨∑-1	н	н	СН3-	Н
A1579	СН3-	\_\^\_\	н	Н	СН3-	Н
A1580	CH3-	\(\rightarrow\)	н	Н	СН3-	н
A1581	CH3-	<u></u>	н	н .	СН3-	н
A1582	CH3-		н	н	СН3-	н
A1583	снз-	<b>○</b> V- <b>〈</b> 〉	н	н	СН3-	н
A1584	СН3-	<u>○</u> V-( <u>&gt;</u> -1	Н	н	СН3-	н
A1585	CH3-	H3CN_N-	Н	н	СН3-	н
A1586	снз-	H3CN N-	н	н	ОН3-	Н

No.	R1	R2	R3	R4	R5	R6
		H3CN N-{_}-				
A1587	CH3-		Н	H	CH3-	Н
A1588	снз–	)' 🖤 '	Н	H	0Н3-	н
A1589	СН3-	, ,	н	н	СН3-	Н
A1590	снз	OCH <sub>3</sub>	н	н	CH3-	Н
A1591	СН3-		н	Н	онз-	Н
A1592	CH3-	CC,	н	н	снз-	Н
A1593	снз-	сн3-	Н	н	снз-	снз-
A1594	CH3-	СН3СН2-	н	н	СН3-	снз
A1595	СН3-	^\\\\ .	Н	н	снз-	снз-
A1596	CH3-	7	Н	н	снз-	СН3-
A1597	CH3-	~~``\	н	н	снз-	СН3
A1598	CH3-	人、	н	н	снз-	СН3
A1599	CH3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	Н	снз-	СН3-
A1600	СН3-	丫	Н	Н	снз-	СН3
A1601	СН3-	· · · · ·	н	Н	СН3-	СН3-
A1602	CH3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	H	Н	снз-	снз-
A1603	снз-	Xx	н	н	СН3-	CH3-
A1604	СН3-	7	н	Н	СН3-	снз-
A1605	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	Н	СН3-	СН3-
A1606	CH3-		н	Н	СН3-	снз-
A1607	СН3-	~~~``\	н	Н	CH3-	снз-
A1608	СН3-	Y~~~	н	н	CH3-	снз-

No.	R1	R2	R3	R4	R5	R6
110.				117	11/0	110
A1609	CH3-	n-C8H17-	н	н	снз-	снз-
A1610	СН3-		н	Н	СН3-	снз-
A1611	СН3-	Q	н	н .	СН3-	снз-
A1612	СН3-		н	н	СН3-	CH3~
A1613	СН3-		н .	н	снз-	CH3-
A1614	снз-	$\triangleright \dashv$	н	н	снз-	СН3-
A1615	СН3-	$\Diamond$	Н	H	снз-	СН3-
A1616	СН3-	$\bigcirc \vdash$	Н	н	снз-	снз-
A1617	СН3-	$\bigcup_{i=1}^{n}$	н	Н	снз-	снз-
A1618	снз-	T	H	Н	снз-	CH3-
A1619	снз-		Н	н	снз-	снз–
A1620	снз-		н	н	снз-	CH3-
A1621	снз-	<u></u>	н	Н	снз-	снз-
A1622	снз-	F →	н	Н	снз-	снз-
A1623	СН3-	<u></u>	н	н	СН3-	снз-
A1624	снз-	F-(	н	н	СН3-	снз-
A1625	СН3-	F-(-)	н	н	снз-	снз-
A1626	СН3-	F———In-{	н	Н	снз-	CH3-
A1627	СН3-	CI C)	н	н	снз-	снз-
A1628	СН3-	CI	н	н	СН3-	СН3-
A1629	СН3-	c⊢ <b>(_</b> )–₁	Н	H ·	снз-	снз-
A1630	CH3-	C <b>├</b>	н	н	СН3-	снз-

No.	R1	R2	R3	R4	R5	. R6
A1631	СН3-	CI-Ond	Н	н	CH3-	СН3-
A1632	снз-	Br	н	Н	СН3-	СН3-
A1633	снз-	Br.	н	Н	CH3-	снз-
A1634	СН3-	Br-{}-{	н	Н	CH3-	· СН3-
A1635	снз-	Br—()	н	н	CH3-	снз-
A1636	снз-	Br-Qm4	н	н	СН3-	снз-
A1637	CH3-		н	н	CH3-	снз-
A1638	СН3-	<u> </u>	н	н	CH3-	снз-
A1639	снз-		н	н	снз-	CH3-
A1640	снз-	CH₃	н	н	CH3-	снз-
A1641	снз-	H <sub>3</sub> C	н	н	СН3-	снз-
A1642	снз-	H <sub>3</sub> C-{\rightarrow}-{\rightarrow}-{\rightarrow}	н	н	CH3-	снз–
A1643	снз-		н	н	. снз-	снз-
A1644	снз-	n-C <sub>3</sub> H <sub>7</sub> -	н	Н	снз-	снз-
A1645	снз-	n-C <sub>4</sub> H <sub>9</sub> -	н	н	снз-	снз-
A1646	снз-	OH →	н	н	снз-	снз-
A1647	СН3-	HO	н	Н	CH3-	снз-
A1648	СН3-	но-{};	н	Н	CH3-	СН3-
A1649	СН3-	OCH <sub>3</sub>	н	н	СН3-	снз-
A1650	СН3-	H₃CO ————————————————————————————————————	н	Н	СН3-	CH3-
A1651	сн3-	H <sub>3</sub> CO-{}-{	н	н	снз-	СН3-
A1652	снз-	H <sub>3</sub> CO-{_}	н	н	CH3-	СН3-

No.	R1	IR2	R3	R4	R5	R6
140	173		170	K4	110	INU
A1653	CH3-	H <sub>3</sub> CO-{}\\\	Н	н	ОН3-	CH3-
A1654	СН3-	OC <sub>2</sub> H <sub>5</sub>	н	н	СН3~	CH3-
A1655	снз-	C <sub>2</sub> H <sub>5</sub> Q	н	н	Сн3-	CH3-
A1656	СН3-	C <sub>2</sub> H <sub>5</sub> O-{_}	н	н	СН3-	CH3-
A1657	снз-	ი-C₃H <sub>7</sub> O-⟨ქ	н	н	СН3-	снз-
A1658	CH3-	n-C₄H <sub>9</sub> O- <b>⟨_</b> _}-{	н	н	СН3-	CH3-
A1659	СН3-	NO <sub>2</sub>	н	н	СН3-	CH3-
A1660	CH3-	O <sub>2</sub> N	н	н	СН3-	СН3-
A1661	CH3-	O <sub>2</sub> N-{	Н	н	СН3-	СН3-
A1662	снз-	CN	н .	Н	СН3-	снз-
A1663	снз-	NC	н	н	СН3-	снз-
A1664	снз-	NC-{}-{	н	н	СН3-	СН3
A1665	онз-	NH <sub>2</sub>	н	Н	СН3-	CH3-
A1666	снз-	H <sub>2</sub> N →	н	н	снз-	СН3-
A1667	снз-	H <sub>2</sub> N-(	н	н	СН3-	СН3-
A1668	СН3-	NMe <sub>2</sub>	н	н	СН3-	CH3-
A1669	снз-	Me₂N	н	н	СН3-	CH3-
A1670	снз-	Me <sub>2</sub> N-(	н	Н	СН3-	СН3-
A1671	снз-		н	Н	СН3-	СН3-
A1672	СН3-	Cn-(	н	Н	СН3-	снз-
A1673	снз-		н	Н	СН3-	снз-
A1674	снз-		н	н	СН3-	СН3-

No.	R1	R2	R3	R4	R5	R6
A1675	СН3		н	н	снз-	CH3-
A1676	СН3-	<u></u>	н	н .	снз-	СН3-
A1677	снз-	<ul><li>~_</li></ul>	Н	H	снз-	СН3-
A1678	СН3-		н	Н	снз-	снз-
A1679	снз-	o_v-⟨>-i	н	Н	снз-	СН3-
A1680	снз-	H3CN N-	н	н	снз-	снз-
A1681	CH3-	H3CN N-⟨}	н	н	снз-	снз-
A1682	снз-		н	Н	снз-	CH3-
A1683	снз-	OCH3	н	н	снз-	СН3
A1684	СН3-	P—COCH <sub>3</sub>	н	н	снз-	CH3-
A1685	CH3-	OCH <sub>3</sub>	н	н	СН3-	снз-
A1686	СН3-	CO .	н	н	снз-	СН3-
A1687	СН3-	OC,	н	Н	СН3-	снз-
A1688	снз-	снз-	н	снз-	снз-	CH3-
A1689	СН3-	снзсн2-	н	снз-	снз-	CH3-
A1690	снз-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	СН3-	снз-	CH3-
A1691	снз-	Y	н	снз	снз-	снз-
A1692	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	СН3-	снз-	СН3-
A1693	СН3-	<u> </u>	н	онз-	онз–	CH3-
A1694	снз-	~~	н	онз-	снз-	СН3-
A1695	СН3-	*	н	CH3-	СН3-	CH3-
A1696	снз-	~~``	н .	снз-	СН3-	CH3-

No.	R1	R2	R3	R4	R5	R6_
A1697	СН3-	Y~~	н	CH3-	снз-	CH3-
A1698	СН3-	人人	н	СН3-	снз-	CH3-
A1699	снз-	个	н	снз-	СН3-	CH3-
A1700	СН3	<b>\\\\</b>	н .	снз-	снз-	CH3-
A1701	СН3-		Н	CH3-	снз-	CH3-
A1702	СН3-	<b>~~</b> ~~``\	н	снз-	СН3-	CH3-
A1703	CH3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	снз-	снз-	CH3-
A1704	CH3-	n-C8H17-	н	снз-	СН3-	CH3-
A1705	CH3-	L~~~	Н	снз-	СН3-	CH3-
A1706	CH3-		Н	СН3-	СН3	CH3-
A1707	CH3-		н	СН3-	СН3-	СН3-
A1708	CH3-		н	CH3-	СН3-	снз-
A1709	CH3-	$\triangleright \rightarrow$	н	CH3-	CH3-	CH3-
A1710	снз-	$\Diamond$ -4	н	СН3-	СН3-	СН3-
A1711	снз		Н	СН3-	СН3-	CH3-
A1712	CH3-	$\bigcirc$	н	СН3-	СН3-	СН3-
A1713	СН3		Н	CH3-	СН3-	CH3-
A1714	СН3-		н	СН3-	CH3-	CH3-
A1715	СН3-		н	CH3-	СН3-	CH3-
A1716	снз–	<b>⊘</b> m{	Н	СН3-	снз-	СН3-
A1717	снз-	F-	н	СН3-	СН3-	снз-
A1718	СН3	<u></u>	н	СН3-	СН3	СН3-

No.	R1	R2	R3	R4	R5	R6
A1719	снз-	F-(-)	н	снз-	снз–	СН3~
A1720	снз-	F-(>-1	Н	СН3-	СН3-	СН3-
A1721	СН3-	F—————————————————————————————————————	н	СН3-	CH3-	СН3-
A1722	снз-	CI →	н	СН3-	СН3-	СН3-
A1723	снз-	CI	н	снз-	снз-	CH3-
A1724	снз-	c⊢(	н	СН3-	СН3-	СН3-
A1725	СН3-	c⊢ <u></u>	н	сн3-	СН3-	СН3-
A1726	снз-	c⊢	н	СН3-	СН3-	снз-
A1727	СН3-	Br	Н	снз-	снз-	CH3
A1728	СН3-	Br. →	н	СН3-	снз-	снз-
A1729	CH3-	Br—{	н	снз-	снз-	CH3-
A1730	СН3-	Br—()	н	СН3-	снз	снз-
A1731	CH3-	Br—Con-{	Н	снз-	снз-	снз-
A1732	CH3-	<b>□</b>	Н	СН3-	снз-	CH3-
A1733	CH3-	 	н	<b>с</b> нз-	онз-	снз-
A1734	ОН3-		н	онз-	снз-	СН3-
A1735	СН3-	CH <sub>3</sub>	н	снз-	снз-	СН3-
A1736	СН3-	H <sub>3</sub> C {	н	СН3-	СН3	СН3-
A1737	СН3-	H <sub>3</sub> C-{_}-{	н	СН3-	снз-	снз-
A1738	снз-		н	СН3-	снз-	снз-
A1739	СН3-	n-C <sub>3</sub> H <sub>7</sub> -{_}-{	н	снз-	снз-	снз-
A1740	снз-	n-C <sub>4</sub> H <sub>9</sub> —{	н	СН3-	снз–	снз-

No.	R1	R2	R3	R4	R5	R6
		OH				
A1741	CH3-		H ———	CH3-	CH3-	CH3-
A1742	СН3-	HO H	н	СН3-	СН3-	снз-
A1743	снз-	но-{>-	н	снз-	СН3-	CH3-
A1744	CH3-	OCH <sub>3</sub>	н	снз-	CH3~	CH3-
A1745	CH3-	H₃CQ	н	снз-	снз-	СН3-
A1746	CH3-	H <sub>3</sub> CO-{}-{	н	снз-	снз-	CH3-
A1747	СН3-	H₃CO- <b>(</b> )	н	снз-	СН3-	СН3~
A1748	снз-	H <sub>3</sub> CO-{	н	снз-	СН3-	CH3-
A1749	снз-	OC <sub>2</sub> H <sub>5</sub>	Н	СН3-	CH3-	CH3-
A1750	СН3-	C <sub>2</sub> H <sub>5</sub> Q	н	снз-	CH3-	СН3-
A1751	СН3-		Н	снз-	СН3-	CH3-
A1752	снз-	n-C <sub>3</sub> H <sub>7</sub> O-	н	СН3-	СН3-	СН3-
A1753	снз-	n-C <sub>4</sub> H <sub>9</sub> O-	Н	СН3-	CH3-	СН3-
A1754	снз-	NO <sub>2</sub>	н	CH3-	СН3-	CH3
A1755	CH3-	O₂N ⟨	н	СН3-	снз-	CH3-
A1756	CH3-	O <sub>2</sub> N-(	Н	СН3-	СН3-	CH3-
A1757	СН3-	CN △	н	сн3-	СН3-	СН3-
A1758	снз-	NC{	н	снз-	СН3-	СН3-
A1759	СН3-	ис-{}-	н	СН3-	СН3-	CH3-
A1760	снз-	NH <sub>2</sub>	Н	снз-	CH3-	снз-
A1761	снз-	H <sub>2</sub> N	н.	СН3-	CH3-	снз-
A1762	снз-	H <sub>2</sub> N-()	н	CH3-	CH3-	СН3-

No.	R1	R2	R3	R4	DE	R6
140.	I I I	NMe <sub>2</sub>	11(0	I T. 4	R5	LZO
A1763	CH3-		н	снз-	СН3-	CH3-
A1764	CH3-	Me <sub>2</sub> N	н	CH3-	СН3~	снз-
A1765	СН3-	Me <sub>2</sub> N-⟨¯⟩∤	н	СН3-	СН3-	СН3
A1766	снз-	CN-\( \( \)	н	CH3-	CH3-	СН3-
A1767	снз-		н	СН3-	CH3~	СН3-
A1768	СН3-		н	СН3-	CH3-	СН3-
A1769	CH3-		н	CH3-	СН3-	CH3-
A1770	СН3-	O1-€}_	н	СН3-	CH3-	СН3-
A1771	CH3-	\\	н	СН3-	СН3-	СН3-
A1772	СН3-		н	СН3	OH3-	СН3-
A1773	СН3-	< <u></u>	Н	CH3-	СН3-	CH3-
A1774	СН3~		н	СН3-	СН3-	СН3-
A1775	СН3-	H³CN_N-∕_>	н	СН3-	СН3-	CH3-
A1776	CH3-	H3CN N-⟨_}	н	снз-	<b>СН3</b> -	СН3-
A1777	СН3-	H3CN N-{_}-{	н	снз-	снз-	CH3-
A1778	СН3-	OCH <sub>3</sub> F-√√	н	снз-	снз <b>–</b>	СН3-
A1779	CH3-	P—C—(	Н	онз-	онз-	CH3-
A1780	СН3-	OCH <sub>3</sub> F-C)···(	н	снз-	СН3-	СН3-
A1781	СН3-	OP	н	снз-	онз-	СН3-
A1782	CH3-		н	CH3-	снз-	СН3-
A1783	снзсн2-	CH3-	н	н	н	н
A1784	снзсн2-	снасн2-	н	н	н	н

	ln4	lpo l		ln.	DE	ne ]
No.	R1	R2	R3	R4	R5	R6
A1785	СН3СН2-		Н	н	н	н
A1786	СН3СН2-	$\nearrow$	н	н	н	Н
A1787	СНЗСН2-	<b>√</b>	н	н	H .	H.
A1788	СН3СН2-	人工	Н	Н	Н	н
A1789	СНЗСН2-	~~`	Н	Н	н	н
A1790	CH3CH2-	Y`	Н	н	Н	н
A1791	СНЗСН2-	<b>^</b> ~\`\	н	Н	н	н
A1792	СНЗСН2-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	н	н
A1793	СНЗСН2-	X	н	Н	н	н
A1794	СНЗСН2-	7	н	н	Н	н
A1795	СНЗСН2-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	Н	н
A1796	СНЗСН2-	L~~	н	Н	н	Н
A1797	СНЗСН2-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	Н	Н	Н
A1798	СНЗСН2-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	н	н
A1799	СНЗСН2-	n-C8H17-	Н	н	н .	н
A1800	СНЗСН2-	\	Н	н	н	Н
A1801	СНЗСН2-		н	н	н	н
A1802	онзон2-		н	н	н	н
A1803	СНЗСН2-		н	Н	Н	н
A1804	CH3CH2-	$\triangleright \rightarrow$	н	н	н	н
A1805	СНЗСН2-	$\Diamond$ -1	н	н	н	н
A1806	СН3СН2-	$\bigcirc$	н	Н	Н	н

No.	R1	R2	R3	R4	R5	R6
A1807	снзсн2-	$\bigcup_{\bot}$	н	н .	Н	н
A1808	снзсн2-	<u></u>	Н	Н	н	Н
A1809	СН3СН2-	<u></u>	н	Н	н	н
A1810	снзсн2-	$\bigcirc$	н	Н	н	н
A1811	снзсн2-		н	н	Н	н
A1812	снзсн2-	<u></u>	Н	н	н	Н
A1813	снзсн2-	F	Н	н	н	Н
A1814	снзсн2-	F-(	н	н	н	H
A1815	СН3СН2-	F-(-)	Н	Н	н	н
A1816	снзсн2-	F-Out	Н	н	н	Н
A1817	СН3СН2-	CI	н	н	Н	н
A1818	СНЗСН2-	CI.	н	Н	Н	Н
A1819	снзсн2-	C├ <b>-</b>	н	н	Н	н
A1820	СНЗСН2-	c⊢ <b>(</b> )–∤	н	н	Н	н
A1821	СНЗСН2-	c⊢©™	Н	н	н	н
A1822	СНЗСН2-	Br	н	н	н .	н
A1823	СН3СН2-	Br.	н	Н	н	н
A1824	СНЗСН2-	Br-{_}-{	н	н	н	н
A1825	СН3СН2-	Br—C>—I	н	н	н	H ·
A1826	СНЗСН2-	Br—∰ייי{	н	н	Н	н
A1827	СНЗСН2-		н	н	н	н
A1828	снзсн2-		Н	Н	Н	н

No.	R1	R2	IR3	R4	R5	R6
				117		
A1829	CH3CH2-		Н	Н	Н	н
A1830	СН3СН2-	CH₃	н	Н .	Н	н
A1831	СНЗСН2-	H₃C <u></u>	н	н	Н	Н
A1832	СНЗСН2-	H <sub>3</sub> C-{}-{	н	Н	н	н
A1833	СНЗСН2-		н	Н	н	H .
A1834	СНЗСН2-	n-C <sub>3</sub> H <sub>7</sub> -{}-{	н	Н	Н	Н
A1835	снзсн2-	n-C <sub>4</sub> H <sub>9</sub> -{}-{	Н	Н	Н	н
A1836	СНЗСН2-	©H ○H	н	Н	н	н
A1837	СНЗСН2-	HO —	н	н	н	Н
A1838	СНЗСН2-	HO-{\bar{\bar{\bar{\bar{\bar{\bar{\bar	Н	н	Н	н
A1839	СНЗСН2-	OCH₃ ◯>⊣	н	Н	Н	Н
A1840	снзсн2-	H <sub>3</sub> CO	н	Н	н	н
A1841	СНЗСН2-	H₃CO-{_}	н	Н	Н	н
A1842	СНЗСН2-	H <sub>3</sub> CO-(	н	Н	н	н
A1843	СНЗСН2-	H3CO-{\_\mathreal_m4	н	Н	н	н
A1844	СНЗСН2-	OC <sub>2</sub> H <sub>5</sub>	Н	н	Н	н
A1845	снзсн2-	C <sub>2</sub> H <sub>5</sub> Q	н	н	н	н
A1846	снзсн2-	C <sub>2</sub> H <sub>5</sub> O-{_}-{	н	н	н	н
A1847	СНЗСН2-	n-C₃H <sub>7</sub> O- <b>⟨</b> _}-	н	н	н	н
A1848	снзсн2-	n-C <sub>4</sub> H <sub>9</sub> O-	н	Н	H	Н
A1849	СН3СН2-	NO <sub>2</sub>	н	н	Н	н
A1850	СНЗСН2-	O <sub>2</sub> N	н	н	н	н

No.	R1	R2	R3	R4	loc .	Inc
140.	1		N3	<del>                                    </del>	R5	R6
A1851	СН3СН2-	O <sub>2</sub> N-{}	Н	н	н	н
A1852	снзсн2-	CN →	Н	н	н	н
A1853	снзсн2-	NC	Н	н	H	н
A1854	снзсн2-	NC-{}	н	Н	н	н
A1855	снзсн2-	NH <sub>2</sub>	Н	н	н	н
A1856	СН3СН2-	H <sub>2</sub> N	Н	н	н	н
A1857	снзсн2-	H <sub>2</sub> N-{}-{	Н	н	н	н
A1858	снзсн2-	NMe <sub>2</sub>	н	н	н	н
A1859	снзсн2-	Me <sub>2</sub> N	н	н	н	н
A1860	снзсн2-	Me <sub>2</sub> N—{	Н	н	H	н
A1861	снзсн2-	Cn-\( \sigma \)	н	H	н	н
A1862	снзсн2-	Cn-<	н	Н	н	н
A1863	снзсн2-	_N-{_}-;	н	н	Н	н
A1864	снзсн2-		н	н	Н	н
A1865	снзсн2-		н	н	Н	н
A1866	снзсн2-		н	Н	н	н
A1867	снзсн2-		н	н	н	н
A1868	снзсн2-		Н	н	н	н
A1869	СНЗСН2-		Н	н	Н	н
A1870	снзсн2-	H³CN N-	н	н.	н	Н
A1871	онзон2-	H₃CN N-	н.	н	Н	Н
A1872	снзсн2-	H₃CN_N-{_}H	Н	н	Н	Н

No.	R1	R2	R3	R4	R5	R6
NO.	1/1	OCH <sub>3</sub>		104		
A1873	CH3CH2-	<b>F</b> -√∑→ ·	н .	H	н	Н
A1874	СН3СН2-	' 🛁 '	Н	Н	Н	н
A1875	снзсн2-	POCH3	Н	н	н	н
A1876	снзсн2-		н	н	н	н
A1877	снзсн2-		Н	н	н	н
A1878	снзсн2-	СН3-	н	СН3-	Н	Н
A1879	СН3СН2-	СН3СН2-	Н	СН3-	н	н
A1880	СН3СН2-	<u> </u>	н	СН3-	н	Н
A1881	СНЗСН2-	$\uparrow \uparrow$	н	СН3-	н	н
A1882	CH3CH2-	<b>~~</b>	н	СН3-	н	н
A1883	онзон2-	人、	н	СН3-	н .	н
A1884	СН3СН2-	$\uparrow \uparrow$	н	снз-	Н	н
A1885	снзсн2-	7	Н	СН3-	Н	Н
A1886	СН3СН2-	~~``	н	снз-	Н	н -
A1887	СН3СН2-	~~	н	снз-	н	н
A1888	СНЗСН2-	Xx	н	СН3-	Н	н
A1889	СНЗСН2-	7	н	снз-	н	н
A1890	СНЗСН2-	<b>\\\\</b>	н	снз-	н	н
A1891	СНЗСН2-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	снз-	н	Н
A1892	снзсн2-	^~^\	н	СН3-	н	Н
A1893	снзсн2-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	СН3-	н	н
A1894	СНЗСН2-	n-C8H17-	н	СН3-	н	н

<u>.</u>						
No.	RI	R2	R3	R4	R5	R6
A1895	снзсн2-		н	СН3-	н	н
A1896	снзсн2-		н	СН3-	Н	н
A1897	СН3СН2-		н	СН3-	H	н
A1898	снзсн2-		н	онз-	н	н
A1899	СН3СН2-	$\triangleright \dashv$	н	СН3-	н	н
A1900	снзсн2-	$\Diamond$	н	CH3-	Н	н
A1901	снзсн2-	$\bigcirc \dashv$	н	СН3-	Н	н
A1902	СНЗСН2-	$\bigcirc$ -4	Н	СН3-	н	н
A1903	CH3CH2-	$\bigcirc$ -i	н	CH3-	Н	Н
A1904	CH3CH2-	<b>△</b> ≀	н	сн3–	н	н
A1905	снзсн2-		н	сн3-	Н	н
A1906	снзсн2-	<b>⊘</b> m4	н	СН3	н	н
A1907	снзсн2-	<b>-</b>	н	сн3–	н	н
A1908	снзсн2-	<b>F</b>	н	снз-	Н	н
A1909	снзсн2-	F-(	н	онз-	н	н
A1910	снзсн2-	F-(>-i	н	снз-	н	Н
A1911	СНЗСН2-	F——>III-{	н	сн3-	Н	н
A1912	снзсн2-	CI →	н	СН3-	н	н
A1913	CH3CH2-	CI	н	СН3-	Н	н
A1914	СНЗОН2-	CI—(	н	сн3-	Н	н
A1915	снзсн2-	C <b>├</b>	н	сн3-	Н	н
A1916	снзсн2-	CI—(	н	онз-	н	н

No.	R1	R2	R3	R4	R5	R6
NO.	IXI	Br			7.0	
A1917	CH3CH2-	<u></u> ;	H	CH3-	Н	Н
A1918	CH3CH2-	Br.	н	СН3-	Н	н .
A1919	CH3CH2-	Br─∰	н	снз-	н	н
A1920	СНЗСН2-	Br—(	Н	снз-	н	Н
A1921	СНЗСН2-	Br—{	н	снз-	н	н
A1922	СНЗСН2-		н	снз-	Н	Н
A1923	СНЗСН2-	<b>\_</b>	н	СН3	н .	н
A1924	СНЗСН2-	H	н	снз-	н	н
A1925	СН3СН2-	CH₃	Н	СН3-	н	н
A1926	снзсн2-	H <sub>3</sub> C{}	н	СН3	н	н
A1927	СНЗСН2-	H <sub>3</sub> C-{}_{{}}{{}	н	СН3-	н	н
A1928	СН3СН2-	C <sub>2</sub> H <sub>5</sub> {	н	СН3-	н	н
A1929	СНЗСН2-	n-C <sub>3</sub> H <sub>7</sub> -{	н	СН3-	н	н
A1930	СНЗСН2-	n-C <sub>4</sub> H <sub>9</sub>	н	СН3-	н	н
A1931	СНЗСН2-	ОН	Н	CH3-	н	н .
A1932	СН3СН2-	HO	н	СН3-	н	н
A1933	СНЗСН2-		н	СН3-	Н	н
A1934	СНЗСН2-	OCH₃	н	СН3-	н	н
A1935	СН3СН2-	H <sub>3</sub> CO	н	СН3	Н	Н
A1936	СН3СН2-	H <sub>3</sub> CO-{_}-{	н	СН3-	Н	н
A1937	СН3СН2-	H₃Cợ- <b>⟨</b> _ <b>&gt;</b> -1	н	СН3-	н	н
A1938	CH3CH2-	H <sub>3</sub> CO-{>\!!!\{	н	CH3-	н	н

No.	R1	R2	R3	loa	loc .	Ina
No.	KI	OC <sub>2</sub> H <sub>5</sub>	143	R4	R5	R6
A1939	CH3CH2-	Co <sub>2</sub> H <sub>5</sub>	н	СН3-	н	н
A1940	CH3CH2-	C <sub>2</sub> H <sub>5</sub> Q	н	СН3-	н	Н
A1941	СН3СН2-	C <sub>2</sub> H <sub>5</sub> O- <b>⟨</b> _){	н	снз-	н	н
A1942	CH3CH2-	n-C <sub>3</sub> H <sub>7</sub> O-	н	СН3-	н	н
A1943	СН3СН2-	n-C <sub>4</sub> H <sub>9</sub> O-	н	снз-	н	н
A1944	СН3СН2-	NO <sub>2</sub>	н	СН3-	н	Н
A1945	СН3СН2-	O <sub>2</sub> N	н	СН3-	н	н
A1946	СН3СН2~	O <sub>2</sub> N-{	н	снз-	н	н
A1947	СН3СН2-	CN CN	н	онз-	н	н
A1948	снзсн2-	NC \	Н	снз-	н	н
A1949	СН3СН2-	NC-{}	н	СН3-	Н	н
A1950	СН3СН2-	NH <sub>2</sub>	н	снз-	Н	н
A1951	СН3СН2-	H <sub>2</sub> N \	н	снз-	н	н
A1952	снзсн2-	H <sub>2</sub> N-{}	н	снз-	Н	н
A1953	СН3СН2-	NMe <sub>2</sub>	н	снз-	н	н
A1954	снзсн2-	Me <sub>2</sub> N	н	снз-	н	н
A1955	СН3СН2-	Me <sub>2</sub> N-{	н	снз–	н	н
A1956	СНЗСН2-		н	снз-	н	н
A1957	СН3СН2-		н	СН3-	н	н
A1958	СНЗСН2-		н	снз-	н	н
A1959	СНЗСН2-		Н	снз-	н .	н
A1960	снзсн2-	O₁-Q	н	снз-	н	н

No.	R1	R2	R3	R4	R5	R6
A1961	СН3СН2-	\(\rightarrow\)	н	СН3-	н	н
A1962	снзсн2-	<b>○</b> / <b>-○</b>	н	СН3-	н .	н
A1963	СНЗСН2-		н	сн3-	Н	н
A1964	СН3СН2-		н	СН3-	н	н
A1965	СНЗСН2-	H3CN_N-	н	CH3-	н	н
A1966	CH3CH2-	H3CN N-	н	CH3-	н	Н
A1967	СНЗСН2-	H3CN N-{_}-{	н	CH3-	н	н
A1968	СНЗСН2-	och,	н	онз-	Н	н
A1969	СНЗСН2-	OCH₃ F—	н	СН3-	Н	н
A197D	СНЗСН2-	OCH3	н	СН3-	н	Н
A1971	СНЗСН2-		н	СН3-	н	Н
A1972	СНЗСН2-	CC,	н	СН3-	н	н

Table-2					
		R <sub>3</sub> R <sub>2</sub> N R <sub>1</sub>			
No	R1	R2	R3_	R4	R5
B1	CH3-	CH3-	Н	н	н .
B2	снз-	СН3СН2-	н	н	н
<b>B</b> 3	CH3-	<b>∕</b> ∕\\	н	н	н
B4	СН3-	74	н	Н	н
B5 <sup>-</sup>	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	Н	н
B6	СН3-	人、	Н	Н	н
<b>B</b> 7	CH3-	丫	н	н	н
B8	СН3-	<b>/</b> ✓ ✓ ✓ ∖	н	н	н
В9	онз-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	н .	н
B10	снз-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н	н	H
B11	снз-	~~~``	н	н	н
B12	СН3-	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	Н	н	н
B13	СН3-		н	н .	н
B14	снз–		Н	н	н
B15	снз-		н	н	н
B16	СН3-		н	Н	Н
B17	снз-	F	Н	н	н
B18	СН3-	F	н	н	, Н

	1_:	· · · · · · · · · · · · · · · · · · ·			
No	R1	R2	R3	R4	R5
B19	СН3-	F-()-{	Ĥ	н	н
B20	СН3-	CI C)	н	н	н
B21	снз-	CI.	н	н	н
B22	СН3-	c⊢(	Н	н	н
B23	СН3-	Br C≯–∤	Н	н	н
B24	снз-	Br.	н	н	н
B25	CH3-	Br—{}	н	Н	н
B26	снз-	CH <sub>3</sub>	н	Н	н
B27	снз-	H <sub>3</sub> C	Н	н	Н
B28	снз-	H₃C- <b>(_</b> )—{	Н	Н	Н
B29	снз-	C <sub>2</sub> H <sub>5</sub> -{}	н	н	н
B30	СН3-	ОН	н	н	н
B31	снз-	но	н	н	н
B32	CH3-	HO-{}-}	н	Н	н
B33	снз-	OCH <sub>3</sub>	н	н	н
B34	СН3-	H <sub>3</sub> CQ	н	н	. н
B35	снз-	H₃CO- <b>(</b> )—{	н	H	н
B36	снз-	C <sub>2</sub> H <sub>5</sub> O-⟨}	н	н	н
B37	снз-	NO <sub>2</sub>	н	н	н
B38	снз-	O <sub>2</sub> N	н	Н	н
B39	снз-	O <sub>2</sub> N-{}	н	н	н

No	RI	R2 CN	R3	R4	R5
B40	снз-		н	H	н
B41	снз-	NC.	н	н	н
B42	СН3-	NC-	н	н	н
B43	СН3-	00	н	н	н
B44	СН3-		н	Н	н
<b>B4</b> 5	CH3-	OD'	н	н	Н
B46	СН3-		Н	н	Н
B47	снз-	PON'	н	н	н
B48	СН3-	Q'i	н	Н	н
B49	СН3-	O day	н	Н	н
B50	СН3-		он	н	н
851	снз-		он	н	н
B52	CH3-	<b>—</b>	он	н	н
B53	снз-	F-()-1	он	н	н
B54	СН3-	CI	он	н	Н
B55	снз-	CI	он	Н	Н
B56	снз-	c⊢ <b>(</b> )→	он	н	Н
<b>B</b> 57	СН3-	Br	он	Н	н
858	снз-	Br.	он	н	н .
B59	СН3-	Br—{}	ОН	Н	н
B60	CH3-	CH₃	он	н	н

1	101	Ino	Inc	754	100
No	R1	R2	R3	R4	R5
B61	CH3-	H <sub>3</sub> C	он	н	Н
B62	CH3-	H <sub>3</sub> C-{	он	Н	н
B63	СН3-	C <sub>2</sub> H <sub>5</sub> —{	он	Н	н
B64	СН3-	ОН	он	Н	н
B65	снз-	HO	ОН	н	H
B66	СН3~	но-{_};	он	н	н
B67	СН3-	OCH <sub>3</sub>	он	н	н
B68	СН3-	H <sub>3</sub> CO	ОН	н	н
B69	СН3-	H <sub>3</sub> CO-	ОН	Н	н
B70	снз-	C <sub>2</sub> H <sub>5</sub> O-{}	он	н	Н
<b>B71</b>	снз-	NO <sub>2</sub>	ОН	Н ,	н
872	СН3-	O <sub>2</sub> N	ОН	н	н
873	снз-	O <sub>2</sub> N-⟨	он	н	Н
B74	снз-	CN	он	н	н
<b>B</b> 75	снз-	NC.	он	н	Н
B76	снз-	NC-{}-{	он	н	н
B77	снз-	aro,	он	Н	н
B78	снз-		он	Н	Н
B79	СН3-	OCC,	он	н .	Н
B80	снз–		CN	. Н	н
B81	CH3	F-I	CN	H ·	. Н

No	RI	Inc			
140	KI	R2	R3	R4	R5
B82	CH3-		CN	н	н
B83	СН3-	F-(	CN	н	. Н
B84	СН3-	CI	CN	н	н
B85	СН3-	CI	CN	н	н
B86	СН3-	c⊢ <u></u>	CN	н .	Н
B87	СН3-	Br	CN	н	н
B88	CH3-	Br.	CN	н	н
B89	снз-	Br—{_}{	CN	Н	Н
B90	снз-	CH₃	CN	Н	Н
B91	CH3-	H <sub>3</sub> C	CN	Н	н
B92	СН3-	H₃C-⟨}	CN	н	н
B93	CH3-		CN	н	н
B94	CH3-	OH	GN	н	н
B95	снз-	HO HO	CN	н	н
B96	снз-	но-{>-;	CN	Н	н
B97	СН3-	OCH₃	CN	н	н
B98	СН3-	H <sub>3</sub> CO	CN	н	н
B99	СН3-	H₃CO-⟨	CN	н	н
B100	СН3-	C <sub>2</sub> H <sub>5</sub> O-{_}-{	GN	Н	н
B101	СН3-	NO <sub>2</sub>	GN	н	н
B102	снз-	O <sub>2</sub> N	CN	н	н

No	R1	R2	R3	R4	R5
140	· · · · ·	4-3	1		1.0
B103	CH3-	O <sub>2</sub> N-{}-{	CN	н	Н
B104	CH3-	CN	GN	Н	н
B105	СН3-	NC	CN	н	н
B106	СН3-	NC-{\rightarrow}-{\lambda}	GN	н	н
B107	CH3-	and,	CN	н	н
B108	СН3-		CN	н	н
B109	CH3-		CN	н	н
B110	CH3-	Н	н	СН3-	н
B111	CH3-	н	Н	СНЗСН2-	н
B112	CH3-	н	н	<b>∼</b> `	Н
B113	CH3-	н	Н	<b>Y</b>	н
B114	СН3-	н	н	~~`	н
B115	CH3-	н	Н	<u></u>	н
B116	СН3	н	н	丫	н
B117	СН3-	Н	н		н
B118	CH3-	н	Н	Y~~	н
B119	СН3-	Н	н	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\	н
B120	СН3-	Н	н		, н
B121	снз-	Н	н	<b>~~~~</b>	\\
B122	GH3-	Н	Н	Qu	н
B123	снз-	н	н	OH OH	н

No	R1	R2	R3	IR4	R5
B124	CH3-	Н	Н	FO	н
B125	CH3-	н	н		н
B126	СН3-	н	н		Н
B127	СН3-	н	н		н
B128	СН3-	н	н	F-1	T.
B129	СН3-	н	н	F.	Н
B130	CH3-	н	н	F-C>-4	Τ
B131	СН3-	н	н	CI CI	н
B132	СН3-	н	н	CI	Н
B133	снз-	н	н	c⊢(	н
B134	СН3-	н	н	CI-	н
B135	снз-	н	н	Br ⊸}	н
B136	снз-	н	н	Br	Н
B137	снз-	н	н	Br—(	н
B138	снз-	н	н	CH₃	н
B139	СН3-	н .	н	H <sub>3</sub> C	н
B140	снз-	н	н	H <sub>3</sub> C-{_}-{	н
B141	снз-	н	н	C <sub>2</sub> H <sub>5</sub>	Н
B142	снз-	н	н	OH OH	н
B143	СН3-	н	н	НО	н
B144	СН3-	н	н	но-(С);	н .

No	R1	R2	100	154	loc
No	IRI		R3	P4 OCH <sub>3</sub>	R5
B145	CH3-	н	Н		н
B146	CH3-	Н	Н	H₃CQ ——;	Н
B147	снз-	Н	Н	H <sub>3</sub> CO- <b>⟨&gt;</b> -{	н
B148	СН3-	н	н	C <sub>2</sub> H <sub>5</sub> O-	Н
B149	CH3-	н	н	NO <sub>2</sub>	Н
B150	снз-	Н	н	O <sub>2</sub> N	н
B151	СН3-	Н	н	O <sub>2</sub> N-(	н
B152	СН3-	н	н	CN →	н
B153	снз–	Н	н	NC.	н
B154	снз-	н	н	NC-()	Н
B155	СН3-	Н	H.		н
B156	снз-	н	н		H
B157	СН3-	Н	Н	F	Н
B158	СН3-	Н	н	FUN	Н
B159	СН3-	н	н	FON	Н
B160	СН3-	Н	н		н
B161	CH3-	Н	Н	\times_N-1	Н
B162	СН3-	н	н	Q\$ <sup>→</sup>	Н
B163	снз-	Н	н	Q <sub>h</sub> ,	н
B164	снз-	н	н	Oh	Н
B165	снз-	н	н	Q <sub>r</sub> ,	н

		· · · · · · · · · · · · · · · · · · ·			
No	R1	R2	R3	R4	R5
B166	снз-	Н	н	CH <sub>3</sub>	н
B167	СН3-	н	Н	H <sup>2</sup> C O	Н
B168	CH3-	н	Н	H <sub>3</sub> C O	н
B169	CH3~	н	Н	<b>◯</b> →	ОН
B170	СН3-	н	н	F	ОН
B171	CH3-	н	Н	F.	он
B172	СН3-	Н	н	F-(	он
B173	СН3-	Н	н	CI	он
B174	СН3-	Н	н	CI	он
B175	СН3-	Н	н	c⊢(	он
B176	CH3-	н	Н	Br	он
B177	СН3-	н	н	Br	он
B178	СН3-	н	н	Br—(i	он
B179	СН3-	н	н	CH <sub>3</sub>	ОН
B180	СН3-	н	Н	H₃C ——;	он
B181	СН3-	н	н	H₃C-⟨{	он
B182	СН3-	Н	н	C <sub>2</sub> H <sub>5</sub> -{_}-{	он
B183	CH3-	н	Н	ОН	ОН
B184	CH3-	н	н	HO	он
B185	снз-	Н	н	но-⟨∑→;	он
B186	CH3-	н	н	OCH₃ C}—{	он

No	R1	R2	R3	R4	R5
			1	H₃CQ	11.0
B187	CH3-	Н	Н		ОН
B188	снз-	Н	H ·	H³CO-{{}	ОН
B189	CH3-	н	н	C <sub>2</sub> H <sub>5</sub> O-	он
B190	CH3-	н	Н	NO <sub>2</sub>	он
B191	CH3-	н	н	O <sub>2</sub> N	он
B192	СН3	н	н	O <sub>2</sub> N-{_}	он
B193	снз-	н	н	CN CN	он
B194	снз-	н	Н	NC	ОН
B195	CH3-	н	н	NC-	он
B196	снз-	Н	Н		ОН
B197	снз-	н	н		он
B198	СН3-	Н	Н		CN
B199	СН3-	н	н	<b>_</b> ;	CN
B200	CH3-	н	н	<u></u> ;	CN
B201	СН3-	Н	Н	F-(	CN
B202	CH3-	н	н	CI	CN
B203	СН3-	н	н	Cı ——	CN
B204	CH3-	Н	н	c⊢ <b>(_</b> }	CN
B205	снз-	Н	Н	Br	CN
B206	Сн3-	н	н	Br.	CN
B207	СН3	н	н	Br <del>-</del> ⟨ <mark>&gt;</mark> -{	СИ

No	IR1	R2	R3	R4	R5
B208	CH3-	н	Н	CH <sub>3</sub>	CN
B209	CH3-	н	н	H₃C —>→	CN
B210	CH3-	Н	н	H₃C-{_}	CN
B211	снз-	н	H.	C <sub>2</sub> H <sub>5</sub> -{}-{	CN
B212	CH3-	н	н	OH ○	CN
B213	СН3-	н	Н	HO	GN
B214	CH3-	н	н	но-(С);	CN
B215	снз-	н	н	OCH <sub>3</sub>	CN
B216	СН3-	Н	Н	H₃CQ —∤	GN
B217	CH3-	н	н	H₃CO-{	CN
B218	СН3-	Н	н	C <sub>2</sub> H <sub>5</sub> O-{}	CN
B219	СН3-	н	н	NO <sub>2</sub>	GN .
B220	СН3-	Н	н	O <sub>2</sub> N	GN
B221	СН3-	н	н	O <sub>2</sub> N-{}_{	CN
B222	CH3~	н	н	CN —	CN
B223	СН3-	Н	н	NC	CN .
B224	СН3-	Н	н	NC-	GN
B225	СН3-	н	н	CQ .	CN
B226	СН3-	н	н	CCC,	CN
B227	СН3-	н	н	<u></u>	000
B228	CH3-	н	н	<b>-</b>	2

No	R1	R2	R3	R4	R5
B229	СН3-	н	Н	F	2
B230	CH3-	н	Н	F-C>-1	<u> </u>
B231	снз-	Н	Н	CI	
B232	ĊH3-	н	н	CI	2,
B233	СН3	н .	н	c⊢ <b>(</b> )→	2
B234	СН3-	Н	н .	Br	<u> </u>
B235	СН3	Н	н	Br.	<u> </u>
B236	CH3	Н	н	Вг-С	1,
B237	СН3-	Н	н	CH₃	<b>}</b> =
B238	СН3-	Н	Н	H₃C —	
B239	CH3-	н	н	H <sub>3</sub> C-\bigcom	
B240	СН3-	Н	н	C <sub>2</sub> H <sub>5</sub> ─{	1,
B241	снз-	н	н	OH OH	<u></u>
B242	СН3-	н	н	HO HO	<u></u>
B243	СН3~	н .	н	но-{}-}	2
B244	СН3-	н	H	OCH <sub>3</sub>	2,
B245	СН3-	Н	Н	H <sub>3</sub> CO	2
B246	CH3-	Н	Н	H3CO-{}-{	2,
B247	СН3-	н	н	C <sub>2</sub> H <sub>5</sub> O-	2
B248	CH3-	н	н	NO <sub>2</sub>	ار کی
B249	СН3-	Н	Н	O <sub>2</sub> N	کہ

No	R1	R2	R3	R4	R5
B250	СН3-	н	Н	O <sub>2</sub> N-\_	°-}
B251	СН3-	н	Н	CN CN	o=\}
B252	CH3-	Н	н	NC T	o=\}
B253	CH3-	Н	н	NC-{}-	o=⟨`
B254	GH3-	Н	Н		o={`
B255	CH3-	н	н		0=

Particularly preferred compounds of the present invention represented by

formula (I) include: 2-(3-Phenylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(4-Fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; 2-(3-(3-Fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; 2-(3-(2-Fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(4-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; (S)-2-(3-(4-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4one: (R)-2-(3-(4-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-2-(3-(3-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Chlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(4-Bromophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(3-Bromophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Bromophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(4-Methylphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(3-Methylphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Methylphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(4-Cyanophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(3-Cyanophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Cyanophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; 2-(3-(4-Methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; 2-(3-(3-Methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one; 2-(3-(2-Methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(2-Ethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; 2-(3-(5-Fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one; 2-(3-(4-Fluoro-3-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one; 2-(3-(4-Fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3Hpyrimidin-4-one; (S)-2-(3-(4-Fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-

pyrimidin-4-one;

- (R)-2-(3-(4-Fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-Chloro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-Fluoro-2-methylphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2-Fluoro-6-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(5-Bromo-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimid in-4-one;
- 2-(3-(2-Bromo-4-fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2-Chloro-6-fluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(2,4-Difluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(2,6-Difluorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(2,6-Dichlorophenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(2,4-Dimethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one:
- 2-(3-(3,4-Dimethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2,5-Dimethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(2,6-Dimethoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(2,4-Difluoro-6-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(5-Cyano-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-Cyano-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;

one:

- 2-(3-(1-Naphthyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one; 2-(3-(2-Naphthyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one; 2-(3-(2,3-Dihydrobenzofuran-7-yl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one; 2-(3-(Benzofuran-2-yl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one; (S)-2-(3-(Benzofuran-2-yl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-
- 2-(3-(4-(Pyrrolidin-1-yl-methyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-(Pyrrolidin-1-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(2-methoxy-4-(pyrrolidin-1-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2-methoxy-5-(pyrrolidin-1-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-(Phenyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one; 2-(3-(4-(4-Fluorophenyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(4-(4-Methoxyphenyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-(2-Methoxyphenyl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(3-(4-(Morpholin-4-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-(4-Methylpiperazin-1-yl)phenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(4-Phenylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(4-Benzylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(4-Benzoylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(4-(1,2-Benzisothiazol-3-yl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(4-Methyl-3-phenylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-Fluoro-2-methoxyphenyl)-4-methylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;

- (S)-2-(3-(4-Fluoro-2-methoxyphenyl)-4-methylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- (R)-2-(3-(4-Fluoro-2-methoxyphenyl)-4-methylpiperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(4-Acetyl-3-(4-fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(4-Benzyl-3-(4-fluoro-2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(4-Benzyl-3-(ethoxycarbonyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(4-methyl-3-(1-naphthyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- 2-(5,5-Dimethyl-3-(2-methoxyphenyl)piperazin-1-yl)-3-methyl-6-(4-pyridyl)-3*H*-pyrimidin-4-one;
- .2-(3-Phenylpiperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-Fluorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(3-Fluorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2-Fluorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-Chlorophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-Bromophenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(4-Methoxyphenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(3-Methoxyphenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- 2-(3-(2-Methoxyphenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- 2-(3-(4-((Pyrrolidin-1-yl)methyl)phenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one;
- (S)-2-(3-(4-(Pyrrolidin-1-yl-methyl)phenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one;
- (R)-2-(3-(4-(Pyrrolidin-1-yl-methyl)phenyl)piperidin-1-yl)-3-methyl-6-(4-pyridyl)-3H-pyrimidin-4-one; and
- 2-(3-Hydroxy-3-phenylpiperidin-1-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one. Salts of the aforementioned preferred compound, and solvates or hydrates of the aforementioned compounds and salts thereof are also preferred.

The 3-substituted-4-pyrimidone compounds represented by the aforementioned formula (I) can be prepared, for example, according to the method

explained below.

(In the above scheme, definitions of R, X and Y are the same as those already described.)

The 2-thiopyrimidone represented by the above formula (III) is prepared easily by a modification of the method described in EP 354,179. The reaction may be carried out in the presence of a base such as sodium hydroxide, potassium hydroxide, sodium methoxide, sodium ethoxide, potassium tert-butoxide, sodium carbonate, sodium hydrogencarbonate, potassium carbonate, triethylamine, diisopropylethylamine, and 1,8-diazabicyclo[5,4,0]undec-7-en for 1 to 100 hours at a suitable temperature ranging from 0 °C to 200 °C under nitrogen or argon atmosphere or under ordinary air to afford the desired compound (III). Examples of a solvent for the reactions include, for example, alcoholic solvent such as methanol, ethanol, 1-propanol, isopropanol, tert-butanol, ethylene glycol, propylene glycol; etheric solvents such as diethyl ether, tert-butyl methyl ether, tetrahydrofuran, isopropyl ether; hydrocarbonic solvents such as benzene, toluene, xylene; halogenated hydrocarbonic solvents such as dichloromethane, chloroform, dichloroethane; aprotic polar solvents such as formamide, N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, dimethyl sulfoxide, sulfolane, hexamethylphosphoric triamide, water and the like. Generally, a single solvent or a mixture of two or more solvents may be used so as to be suitable to a base used.

Then the 2-thiopyrimidone derivative (III) is transformed into the 2-chloropyrimidone (IV) by a chlorinating agent. The reaction time and temperature depend on the chlorinating agent used. Examples of a chlorinating agent for the reactions include, for example, thionyl chloride, thionyl chloride and dimethylformamide, phosphorus oxychloride, phosphorus oxychloride and

dimethylformamide, oxalyl chloride, phosphorous oxychloride and dimethylformamide, and phosphorus pentachloride.

The amine represented by the above formula (V) may be prepared by a modification of the method described in Japanese Patent Unexamined Publication [Kokai] No. 52-139085/1977 or according to well-known methods of one skilled in the art.

Then the chloride derivative (IV) is allowed to react with the amine (V) or salts thereof in the presence of a base such as sodium hydroxide, potassium hydroxide, sodium methoxide, sodium ethoxide, sodium carbonate, sodium hydrogencarbonate, potassium carbonate, triethylamine, diisopropylethylamine, and 1,8-diazabicyclo[5,4,0]undec-7-en for 0.1 to 100 hours at a suitable temperature ranging from 0 °C to 200 °C under nitrogen or argon atmosphere or under ordinary air to afford the desired compound (II).

Examples of a solvent for the reactions include, for example, alcoholic solvent such as methanol, ethanol, 1-propanol, isopropanol, tert-butanol, ethylene glycol, propylene glycol; etheric solvents such as diethyl ether, tert-butyl methyl ether, tetrahydrofuran, isopropyl ether; hydrocarbonic solvents such as benzene, toluene, xylene; halogenated hydrocarbonic solvents such as dichloromethane, chloroform, dichloroethane; aprotic polar solvents such as formamide, N,N-dimethylformamide, N,N-dimethylacetamide, N-methylpyrrolidone, dimethyl sulfoxide, sulfolane, hexamethylphosphoric triamide, water and the like. Generally, a single solvent or a mixture of two or more solvents may be used so as to be suitable to a base used.

The compounds of the present invention have inhibitory activity against TPK1, and they inhibit TPK1 activity in neurodegenerative diseases like Alzheimer disease, thereby suppress the neurotoxicity of A $\beta$  and the formation of PHF and inhibit the nerve cell death. Accordingly, the compounds of the present invention are useful as an active ingredient of a medicament which radically enables preventive and/or therapeutic treatment of Alzheimer disease. In addition, the compounds of the present invention are also useful as an active ingredient of a medicament for preventive and/or therapeutic treatment of ischemic cerebrovascular accidents, Down syndrome, cerebral bleeding due to solitary

cerebral amyloid angiopathy, progressive supranuclear palsy, subacute sclerosing panencephalitis, postencephalitic parkinsonism, pugilistic encephalosis, Guam parkinsonism-dementia complex, Lewy body disease, Pick's disease, corticobasal degeneration frontotemporal dementia, vascular dementia, acute stroke and traumatic injuries, brain and spinal cord trauma, peripheral neuropathies, retinopathies and glaucoma, non-insulin dependent diabetes (such as diabetes type II), and obesity, manic depressive illness, schizophrenia, alopecia, cancers such as breast cancer, non-small cell lung carcinoma, thyroid cancer, T or B-cell leukemia and several virus-induced tumors.

As the active ingredient of the medicament of the present invention, a substance may be used which is selected from the group consisting of the compound represented by the aforementioned formula (I) and pharmacologically acceptable salts thereof, and solvates thereof and hydrates thereof. The substance, per se, may be administered as the medicament of the present invention, however, it is desirable to administer the medicament in a form of a pharmaceutical composition which comprises the aforementioned substance as an active ingredient and one or more of pharmaceutical additives. As the active ingredient of the medicament of the present invention, two or more of the aforementioned substance may be used in combination. The above pharmaceutical composition may be supplemented with an active ingredient of other medicament for the treatment of, for example, Alzheimer disease, vascular dementia, acute stroke and traumatic injuries, brain and spinal cord trauma, peripheral neuropathies, retinopathies and glaucoma, non-insulin dependent diabetes (such as diabetes type II), and obesity, manic depressive illness, schizophrenia, alopecia, cancers such as breast cancer, non-small cell lung carcinoma, thyroid cancer, T or B-cell leukemia and several virus-induced tumors.

A type of the pharmaceutical composition is not particularly limited, and the composition may be provided as any formulation for oral or parenteral administration. For example, the pharmaceutical composition may be formulated, for example, in the form of pharmaceutical compositions for oral administration such as granules, fine granules, powders, hard capsules, soft capsules, syrups, emulsions, suspensions, solutions and the like, or in the form of pharmaceutical compositions for parenteral administrations such as injections for intravenous,

intramuscular, or subcutaneous administration, drip infusions, transdermal preparations, transmucosal preparations, nasal drops, inhalants, suppositories and the like. Injections or drip infusions may be prepared as powdery preparations such as in the form of lyophilized preparations, and may be used by dissolving just before use in an appropriate aqueous medium such as physiological saline. Sustained-release preparations such as those coated with a polymer may be directly

administered intracerebrally.

Types of pharmaceutical additives used for the manufacture of the pharmaceutical composition, content rations of the pharmaceutical additives relative to the active ingredient, and methods for preparing the pharmaceutical composition may be appropriately chosen by those skilled in the art. Inorganic or organic substances, or solid or liquid substances may be used as pharmaceutical additives. Generally, the pharmaceutical additives may be incorporated in a ratio ranging from 1% by weight to 90% by weight based on the weight of an active ingredient.

Examples of excipients used for the preparation of solid pharmaceutical compositions include, for example, lactose, sucrose, starch, talc, cellulose, dextrin, kaolin, calcium carbonate and the like. For the preparation of liquid compositions for oral administration, a conventional inert diluent such as water or a vegetable oil may be used. The liquid composition may contain, in addition to the inert diluent, auxiliaries such as moistening agents, suspension aids, sweeteners, aromatics, colorants, and preservatives. The liquid composition may be filled in capsules made of an absorbable material such as gelatin. Examples of solvents or suspension mediums used for the preparation of compositions for parenteral administration, e.g. injections, suppositories, include water, propylene glycol, polyethylene glycol, benzyl alcohol, ethyl oleate, lecithin and the like. Examples of base materials used for suppositories include, for example, cacao butter, emulsified cacao butter, lauric lipid, witepsol.

Dose and frequency of administration of the medicament of the present invention are not particularly limited, and they may be appropriately chosen depending on conditions such as a purpose of preventive and/or therapeutic treatment, a type of a disease, the body weight or age of a patient, severity of a disease and the like. Generally, a daily dose for oral administration to an adult may be 0.01 to 1,000 mg (the weight of an active ingredient), and the dose may be administered once a day or several times a day as divided portions, or once in several days. When the medicament is used as an injection, administrations may preferably be performed continuously or intermittently in a daily dose of 0.001 to 100 mg (the weight of an active ingredient) to an adult.

## Examples

The present invention will be explained more specifically with reference to examples. However, the scope of the present invention is not limited to the following examples. The compound numbers in the examples correspond to those in the table above.

Reference Example 1: Synthesis of 2-mercapto-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one

A solution of ethyl 3-oxo-3-(4-pyridyl)propionate (29.0 g, 150 mmol), N-methyl thiourea (40.6 g, 450 mmol) and 1,8-diazabicyclo[5,4,0]-7-undecene (22.4 ml, 150 mmol) was refluxed for 4 hours and the solution of methanesulfonic acid (14.4 g, 150 mmol) in water (50 ml) was added after cooling by ice-water. The precipitate was washed with water, filtered and dried to give the title compound (23.7 g, 72%).

<sup>1</sup>H-NMR (DMSO-d<sub>6</sub>)  $\delta$ : 3.58(s, 3H), 6.40(s, 1H), 7.72(dd, J=1.8, 4.5Hz, 2H), 8.73(dd, J=1.5, 4.8Hz, 2H), 12.92(brd, 1H).

Reference Example 2: Synthesis of 2-chloro-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one

Phosphorous oxychloride (26.11g, 170 mmol) was added to dimethylformamide(180 ml) and stirred 20 min. 2-Mercapto-3-methyl-6-(4-pyridyl)-pyrimidine-4-one (24.15 g, 110 mmol) was added to the solution and stirred 5 min and then stirred at 70°C for 2 hours. Ethyl acetate (630 ml) was added to the ice-cooled solution and precipitate was collected by filtration after stirring for 20 minutes. After drying, the precipitate was dissolved in water (400 ml) and pH was

adjusted to 10 by using aqueous sodium hydroxide. The precipitate was washed with water, filtered and dried to give the title compound (18.82 g, 77%).  $^{1}$ H-NMR (CDCl<sub>3</sub>)  $\delta$ : 3.72(s, 3H), 6.90(s, 1H), 7.78(dd, J=1.7, 4.5Hz, 2H), 8.75(dd, J=1.6, 4.5Hz, 2H).

Example 1: Synthesis of 2-(2-(4-fluoro-2-methoxyphenyl)piperazin-4-yl)-3-methyl-6-(4-pyridyl)-3 H-pyrimidin-4-one dihydrochloride

A solution of 2-bromo-5-fluoroanisole (11.8 g, 57.6 mmol) in tetrahydrofuran (60 ml) was dropped into the magnesium (1.40 g, 57.6 mmol) in refluxed tetrahydrofuran (32 ml) containing small amount of 1,2-dibromoethane and refluxed for 15 min. After addition of tetrehydrofuran (50 ml), the solution was cooled to -78 °C and diethyl oxalate (7.41 g, 50.7 mmol) in diethyl ether (50 ml) was dropped into the solution. After stirring at same temperature for 30 min, the solution was warmed to -10°C and 1N aqueous hydrogen chloride (50 ml) and water were added. Organic layer was extracted with diethyl ether, washed with brine and dried over magnesium sulfate. After removal of the solvent under reduced pressure, purification of the residue by silica gel column chromatography (eluent: hexane/ethyl acetate = 5/2) gave ethyl 2-(4-fluoro-2-methoxyphenyl)-2-oxoacetate (6.80g, 59%)

<sup>1</sup>H-NMR (CDCl<sub>3</sub>)  $\delta$ : 1.40(3H, t, J=7.1 Hz),3.87(3H, s), 4.89(2H, q, J=7.1Hz), 6.68(1H, d, J=10.5 Hz), 6.77-6.81(1H, m), 7.91-7.95(1H, m).

Ethylenediamine (0.60 g, 10.0 mmol) was added to a solution of ethyl 2-(4-fluoro-2-methoxyphenyl)-2-oxoacetate (2.26 g, 10.0 mmol) in ethanol(30 ml) and refluxed 4 hr. After removal of the solvent under reduced pressure, residue was washed with ethanol-diethyl ether to give 5,6-dihydro-3-(4-fluoro-2-methoxyphenyl)pyrazinone (1.76 g, 79%).

<sup>1</sup>H-NMR (CDCl<sub>3</sub>)  $\delta$ : 3.50-3.56 (2H, m), 3.81 (3H, s), 3.88-3.92 (2H, m), 6.65(1H, d, J=11.0 Hz), 6.70-6.76 (1H, m), 6.89(1H, bs), 7.36-7.40(1H, m).

5,6-Dihydro-3-(4-fluoro-2-methoxyphenyl)pyrazinone was added to the solution of lithium aluminium hydride (0.46 g, 12 mmol) in diethyl ether (25 ml) and refluxed for 6 hr. Water (0.48 ml), 15% sodium hydroxide solution (0.48 ml) and again water (1.21 ml) were added to the ice-cooled solution and the precipitate was

filtered and washed with dichloromethane. Combined organic layer was evaporated to give 2-(4-fluoro-2-methoxyphenyl)piperazine (0.83 g, 99%).

<sup>1</sup>H-NMR (CDCl<sub>3</sub>)  $\delta$ : 2.02(2H, s), 2.57-2.63 (1H, m), 2.80-2.89 (1H, m), 2.92-2.99 (2H, m), 3.06-3.12 (2H, m), 3.80(3H, s), 4.06 (1H, d, J=10.0 Hz), 6.56-6.65 (2H, m), 7.40 (1H, t, J=7.8 Hz).

2-Chloro-3-methyl-6-(4-pyridyl)-pyrimidin-4-one (222 mg, 1.0 mmol) was added to an ice-cooled solution of 2-(4-fluoro-2-methoxyphenyl)piperazine (210 mg, 1.0 mmol), triethylamine (0.15 ml, 1.1 mmol) in N,N-dimethylformamide (10 ml) and stirred at that temperature for 1 hr and then at room temperature for 2 hr. Next day, reaction was quenched by ice-water and the filtrate was washed with water and dried to give 2-(2-(4-fluoro-2-methoxyphenyl)piperazin-4-yl)-3-methyl-6-(4-pyridyl)-pyrimidin-4-one (246 mg, 62%).

<sup>1</sup>H-NMR (CDCl<sub>3</sub>)  $\delta$ : 2.89-2.96 (1H, m), 3.19-3.31 (3H, m), 3.59 (3H, s), 3.62-3.74 (2H, m), 3.85 (3H, s), 4.39-4.44 (1H, m), 6.63-6.71 (2H, m), 6.67 (1H, s), 7.51-7.55 (1H, m), 7.81 (2H, dd, J=1.7, 4.6 Hz), 8.71 (2H, dd, J=1.7, 4.6 Hz).

4N Hydrogen chloride in 1,4-dioxane (0.4 ml) was added to the solution of 2-(2-(4-fluoro-2-methoxyphenyl)piperazin-4-yl)-3-methyl-6-(4-pyridyl)-pyrimidin-4-one (217 mg, 0.6 mmol) in dichloromethane (5 ml) and stirred for 15 min. After addition of diethyl ether, filtration and wash with diethyl ether and dryness gave the title compound (260 mg, quant.).

Example 2: Synthesis of 2-(2-(4-methoxyphenyl)-piperazine-4-yl)-3-methyl-6-(4-pyridyl)pyrimidin-4-one dihydrochloride

Dimethylslufoxide (50 ml) solution of 4-methyoxyphenacylbromide (9.94 g, 43.4 mmol) and water (1.6 ml, 88.8 mmol) were stirred at 50°C for 2.5 hr. Water was added and the solution was extracted with ethyl acetate 3 times and washed with brine and then dried over sodium sulfate. Removal of the solvent gave 4-methoxyphenylglyoxal (8.30 g, quant.).

<sup>1</sup>H-NMR (DMSO)  $\delta$ : 3.84 (3H, s), 6.60-6.69 (1H, m), 7.04 (2H, d, J=8.8 Hz), 8.05 (2H, d, J=9.1 Hz).

Methanol (5 ml) solution of ethylenediamine (3.74 g, 62.29 mmol) was added to the ice-cooled solution of 4-methoxyphenylglyoxal (8.30 g, 45.5 mmol) in

methanol (100 ml) and tetrahydrofuran (50 ml) and stirred for 10 min. After cooling to 0°C, sodium tetrahydroborate (6.14 g, 162.2 mmol) and additional methanol (50 ml) was added and stirred overnight. After removal of the solvent, aqueous sodium hydroxide was added and was extracted with dichloromethane three times and washed with brine and dried over sodium sulfate. After removal of the solvent, purification of the residue by silica gel column chromatography (eluent; dichloromethane/ethanol/diethylamine = 20/2/1) gave 2-(4-methoxypheny)-piperazine (3.96 g, 45%).

<sup>1</sup>H-NMR (CDCl<sub>3</sub>)  $\delta$ : 2.69(1H, dd, J=10.3, 11.9 Hz), 2.80-3.01(4H, m), 3.07-3.11 (1H, m), 3.68-3.73(1H, m), 3.79(3H, s), 6.84-6.88 (2H, m), 7.27-7.32 (2H, m).

A solution of triethylamine (697 mg, 6.9 mmol), 2-(4-methoxyphenyl)-piperazine (430 mg, tetrahydrofuran (10 ml) was stirred at room temperature for 30 min and at 50°C for 3 hr. Solvent was removed under reduced pressure, and 1N aqueous sodium hydroxide solution was added to the residue and extracted by dichloromethane three times and washed with brine and dried over sodium sulfate. After removal of the solvent under reduced pressure, the residue was purified by silica gel column chromatography (eluent; dichloromethane/ethanol = 10/1) to give 2-(2-(4-methoxyphenyl)-piperazine-4-yl)-3-methyl-6-(4-pyridyl)pyrimidin-4-one (594 mg, 76%)

<sup>1</sup>H-NMR (CDCl<sub>3</sub>)  $\delta$ : 3.02 (1H, dd, J=10.8, 12.7 Hz), 3.18-3.25 (3H, m), 3.55 (3H, s), 3.57-3.67 (2H, m), 3.82 (3H, s), 3.98(1H, dd, J=2.7, 10.8 Hz), 6.67 (1H, s), 6.92 (2H, d, J=8.7 Hz), 7.37 (2H, d, J=8.7 Hz), 7.80 (2H, d, J=6.0 Hz), 8.71 (2H, d, J=6.0 Hz).

4N Hydrogen chloride in ethyl acetate (5 ml) was added to the solution of 2-(2-(4-methoxyphenyl)-piperazine-4-yl)-3-methyl-6-(4-pyridyl)pyrimidin-4-one (594 mg, 1.6 mmol) in dichloromethane (5 ml) and stirred for 1 hr. Wash with ethyl acetate after removal of the solvent and dryness gave the title compound (683 mg, 96%).

Example 3: Synthesis of 2-(2-(4-chlorophenyl)-piperazine-4-yl)-3-methyl-6-(4-pyridyl)pyrimidin-4-one hydrochloride

Mixture of methyl (4-chlorophenyl)acetate (5.10 g, 27.6 mmol) and N-bromosuccinimide (5.16 g, 29 mmol) in carbon tetrachloride was treated by Hg

lamp. After filtration, solvent was removed under reduced pressure and the residue was dissolved in methanol. Ethylenediamine (2.03 ml, 30.4 mmol) and triethylamine (2.06 ml, 14.8 mmol) and di-tert-butyldicarbonate (3.10 ml, 13.5 mmol) were added to the solution of 3-(4-chlorophenyl)piperazin-2-one (2.60 g, 12.3 mmol) in dichloromethane (100 ml) and stirred. The reaction mixture was washed with 1N aqueous hydrogen chloride, water, brine and then dried. After removal of the solvent under reduced pressure, residue was purified by silica gel column chromatography to give 4-(tert-butoxycarbonyl)-3-(4-chlorophenyl)-piperazin-2-one.

<sup>1</sup>H-NMR (CDCl<sub>3</sub>)  $\delta$ : 1.44 (9H, s), 3.21-3.32 (2H, m), 3.48 (1H, m), 4.04 (1H, brs), 5.66 (1H, brs), 7.10 (1H, brs), 7.30-7.38 (4H, m).

Solution of 4-(tert-butoxycarbonyl)-3-(4-chlorophenyl)-piperazin-2-one (500 mg, 1.6 mmol) and acetic acid (929 μl, 16 mmol) were added to a refluxed solution of sodium borohydride (608 mg, 16 mmol) in 1,4-dioxane (5 ml) and reflux was continued. The reaction was quenched by water and extracted with dichloromethane and washed with brine and dried. After removal of the solvent, residue was purified by silica gel column chromatography to give 4-(tert-butoxycarbonyl)-3-(4-chlorophenyl)piperazine (330 mg, 69%).

1H-NMR (CDCl<sub>3</sub>) δ: 1.46(9H, s), 2.76-2.99(3H, m), 3.13(1H, dd, J=13.0 Hz, 4.3 Hz), 3.45-3.49(2H, m), 3.92(1H, m), 5.15(1H, s), 7.27-7.33(4H, m).

A solution of 4-(tert-butoxycarbonyl)-3-(4-chlorophenyl)piperazine (330 mg, 1.1 mmol), 2-chloro-3-methyl-6-(4-pyridyl)pyrimidin-4-one (246 mg, 1.1 mmol) and triethylamine (170  $\mu$ l, 1.22 mmol) in tetrahydrofuran were refluxed. Usual workup and purification by silica gel column chromatography gave 2-(1-(tert-butoxy-carbonyl)-2-(4-chlorophenyl)-piperazine-4-yl)-3-methyl-6-(4-pyridyl)pyrimidin-4-one (500 mg, 93%).

<sup>1</sup>H-NMR (CDCl<sub>3</sub>)  $\delta$ : 1.45(9H, s), 3.09(1H, m), 3,35(3H, s), 3.40-3.63(4H, m), 3.96-4.19(2H, m), 5.43(1H, s), 6.68(1H, s), 7.23(2H, d, J=8.3 Hz), 7.32(2H, d, J=8.3 Hz), 7.78(2H, d, J=5.9 Hz), 8.72(2H, d, J=5.9 Hz).

4N Hydrogen chloride in ethyl acetate was added to the solution of 2-(1-(tert-butoxycarbonyl)-2-(4-chlorophenyl)-piperazine-4-yl)-3-methyl-6-(4-pyridyl)pyrimidin-4-one (500 mg, 1.0 mmol) in ethyl acetate and stirred. Filtration

and successive dryness gave the title compound (373mg, 79%).

Example 4: Synthesis of 3-methyl-2-(3-(4-1-pyrrolidinyl)methylphenyl)piperidine-1-yl)-6-(4-pyridyl)pyrimidin-4-one fumarate

Tetrakis(triphenylphosphine)palladium (0.65 g, 0.56 mmol),
4-formylphenylboric acid (2.81 g, 18.7 mmol), 2M aqueous sodium carbonate (18.7 ml, 37.4 mmol) and ethanol were added to the nitrogen-saturated solution of
3-bromopyridine (2.66 g, 16.8 mmol) in toluene and refluxed under nitrogen for 8 hrs. Water was added to the solution and extracted with ethyl acetate, washed with water and brine and dried. Solvents were removed under reduced pressure and the residue was purified by silica gel column chromatography (eluent; hexane/ethyl acetate = 1/1.5) to give 4-(3-pyridyl)benzaldehyde (0.78 g, 25%).

Methyl iodide (0.8 ml, 12.9 mmol) was added to a solution of 4-(3-pyridyl)benzaldehyde (0.78 g, 4.3 mmol) in dichloromethane and stirred 2 days. Additional methyl iodide (0.8 ml, 12.9 mmol) was added and stirred for 3 hr. After removal of the solvent, methanol was added to the residue and ice-cooled. Sodium tetrahydroborate (6.4 g, 17.0 mmol) was added to the solution and stirred for 1.5 hr with warming to room temperature. Organic solvents were removed under reduced pressure after addition of water and extracted with ethyl acetate, washed with water and brine and dried over sodium sulfate. After removal of the solvent under reduced pressure, residue was purified by silica gel chromatography (eluent ethyl acetate to methanol) to give 3-(4-hydroxymethylphenyl)-1-methyl-1,2,5,6-tetrahydropyridine (0.63 g, 72%).

Triethylamine (1.29 ml, 9.2 mmol), acetic anhydride (0.35 ml, 3.7 mmol) were added to a solution of 4-(hydroxymethyl)phenyl-1-methyl-1,2,5,6-tetrahydropyridine (0.63 g, 3.1 mmol) in dichloromethane and stirred overnight. Organic solvents were removed under reduced pressure after addition of water and extracted with ethyl acetate, washed with water and brine and dried over sodium sulfate. Removal of the solvent under reduced pressure gave 3-(4-acetozymethyl-phenyl)-1-methyl-1,2,5,6-tetrahydropyridine (0.67 g, 89%).

A solution of 3-(4-acetoxymethylphenyl)-1-methyl-1,2,5,6tetrahydropyridine (0.67 g, 2.7 mmol) and 1-chloroethyl chloroformate (0.36 ml, 3.3 mmol) in dichloroethane was refluxed for 2 hr. Organic solvents were removed under reduced pressure after addition of water and extracted with ethyl acetate, washed with water and brine and dried over sodium sulfate. After removal of the solvent, methanol was added and refluxed for 1.5 hr. Tetrahydrofuran and water were added to the residue after removal of the solvent under reduced pressure and triethylamine (1.9 ml, 13.6 mmol) and di-tert-butyl dicarbonate (0.66 g, 3.0 mmol) were added and stirred. Organic solvents were removed under reduced pressure after addition of water and extracted with ethyl acetate, washed with water and brine and dried over sodium sulfate. Removal of the solvent under reduced pressure and the residue was purified by silica gel chromatography to give 3-(4-acetoxymethylphenyl)-1-(tert-butoxycarbonyl)-1,2,5,6-tetrahydropyridine (0.71 g, 78%).

Palladium on charcoal was added to the solution of 3-(4-acetoxymethylphenyl)-1-(tert-butoxycarbonyl)-1,2,5,6-tetrahydropyridine (0.71 g, 2.1 mmol) in ethyl acetate and stirred under hydrogen atmosphere. After filtration with celite and removal of the solvent under reduced pressure, methanol and 1N aqueous sodium hydroxide were added and stirred. Organic solvents were removed under reduced pressure after addition of water and extracted with ethyl acetate, washed with water and brine and dried over sodium sulfate. Removal of the solvent under reduced pressure and the residue was purified by silica gel chromatography (eluent; hexane/ethyl acetate = 3/1) to give 3-(4-hydroxymethylphenyl)-1-(tert-butoxycarbonyl)piperidine (0.39 g, 62%).

Triethylamine (0.47 g, 3.4 mmol) and methanesulfonyl chloride (0.12 ml, 1.6 mmol) were added to an ice-cooled solution of 3-(4-hydroxymethylphenyl)-1-(tert-butoxycarbonyl)piperidine (0.39 g, 1.34 mmol) in dichloromethane and stirred for 7.5 hr. Pyrrolidine (1.0 ml, 12 mmol) was added to the solution and stirred overnight. Organic solvents were removed under reduced pressure after addition of water and extracted with ethyl acetate, washed with water and brine and dried over sodium sulfate. Removal of the solvent under reduced pressure and the residue was purified by silica gel chromatography (eluent; ethyl acetate to ethyl acetate/methanol = 1/1, then methanol only) to give 3-(4-(1-pyrrolidinyl)methyl-phenyl)-1-(tert-butoxycarbonyl)piperidine (0.26 g, 56%).

4N Hydrogen chloride in ethyl acetate was added to 3-(4-(1-pyrrolidinyl)-methylphenyl)-1-(tert-butoxycarbonyl)piperidine (0.26 g, 0.75 mmol) and stirred overnight. After filtration and dryness, triethylamine (0.5 ml, 3.6 mmol), 2-chloro-3-methyl-6-(4-pyridyl)-pyrimidin-4-one (0.14 g, 0.63 mmol) and tetrahydrofuran were added and stirred at 70°C. Organic solvents were removed under reduced pressure after addition of water and extracted with ethyl acetate, washed with water and brine and dried over sodium sulfate. Removal of the solvent under reduced pressure and the residue was dissolved into ethyl acetate. A solution of fumaric acid (0.095 g, 0.82 mmol) in acetone was added and the resulting precipitate was filtered and dried to give the title compound (0.29 g, 76%).

Example 5: Synthesis of (R)-2-(2-(4-chlorophenyl)piperazin-4-yl)-3-methyl-6-(4-pyridyl)-pyrimidin-4- one

To a solution of (S)-2-methyl-CBS-oxazaborolidine (27.6 mL, 1.0 M solution in toluene, 27.6 mmol) was added borane-tetrahydrofuran complex (166 ml, 1.0 M solution in tetrahydrofuran, 166 mmol) at -40 °C. To the resulting solution was added a solution of 4'-chlorophenacyl bromide (32.25 g, 138.1 mmol) in tetrahydrofuran (200 ml) through dropping funnel over 1 h at -40 °C. After stirring for 3 hours below 0 °C, methanol (ca. 50 ml) was added dropwise. After stirring the resulting solution for additional 30 min at room temperature, solvent was removed under reduced pressure. The residue, dissolved in ethyl acetate, was treated with 1 N hydrochloric acid to form white precipitate, which was filtered off. The layers of the filtrate was separated, and the organic layer was washed with hydrochloric acid and brine successively, dried over anhydrous sodium sulfate, filtered, and concentrated under reduced pressure. The residue was used for the next reaction without further purification.

The residue was dissolved in ether (250 ml), and stirred with potassium hydroxide (15.5 g, 276 mmol) in water (250 ml) vigorously. After consumption of the starting material, the layers were separated. The organic layer was washed with brine, dried over anhydrous sodium sulfate, filtered, and concentrated under reduced pressure. The residue was used for the next reaction without further purification.

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The residue was heated with benzylamine (37.7 ml, 345 mmol) at 80 °C for 4.5 h. After cooling at room temperature, the resulting white crystals was washed with ether/hexane and collected to afford (S)-2-benzylamino-1-(4-chlorophenyl)-ethanol (23.8 g, 65.8%). The excess benzylamine in the filtrate was distilled off at 120 °C under reduced pressure. From the residue, another (S)-2-benzylamino-1-(4-chlorophenyl)ethanol (2.41 g, 6.7%) was obtained.

¹H NMR (CDCl<sub>3</sub>) ™: 2.68(1H, dd, J=12.3, 8.9Hz), 2.92(1H, dd, J=12.3, 3.7Hz), 3.80(1H, d, J=11.9Hz), 3.86(1H, d, J=11.9Hz), 4.68(1H, dd, J=8.9, 3.7Hz), 7.30(9H, m).

To a suspension of (S)-2-benzylamino-1-(4-chlorophenyl)ethanol (15.76 g, 60.21 mmol) and triethylamine (33.6 ml, 241 mmol) in dichloromethane (300 ml) was added a solution of thionyl chloride (4.83 ml, 66.2 mmol) in dichloromethane (20 ml) at -78 °C over 20 min. The resulting suspension was stirred at -78 °C for 20 min and at 0 °C for additional 20 min. The reaction mixture was partitioned between ether and water, and the organic layer was washed with brine, dried over anhydrous sodium sulfate, filtered, and concentrated under reduced pressure. The residue was purified by silica gel column chromatography (eluent: 10-20% ethyl acetate-hexane) to afford (2RS,5S)-3-benzyl-5-(4-chlorophenyl)-1,2,3-oxathiazolidine 2-oxide (16.2 g 87.4%) as a pale yellow solid.

The resulting product was obtained as a mixture of two diastereomers due to the S-oxide.

major isomer:  ${}^{1}$ H\_NMR (CDCl<sub>8</sub>)  $\delta$ : 3.31(1H, dd, J=10.5, 9.9Hz), 3.55(1H, dd, J=9.0, 6.3Hz), 3.88(1H, d, J=13.2Hz), 4.37(1H, d, J=13.2Hz), 5.49(1H, dd, J=10.5, 6.3Hz), 7.22-7.43(9H, m).

minor isomer: <sup>1</sup>H NMR (CDCl<sub>8</sub>)  $\delta$ : 3.21(1H, dd, J=13.5, 4.5Hz), 3.77(1H, dd, J=13.5, 11.4Hz), 4.05(1H, d, J=13.5Hz), 4.38(1H, d, J=13.5Hz), 5.99(1H, dd, J=11.4, 4.5Hz), 7.22-7.43(9H, m).

A solution of (2RS,5S)-3-benzyl-5-(4-chlorophenyl)-1,2,3-oxathiazolidine 2-oxide (16.2 g, 52.6 mmol) and sodium azide (17.11 g, 263.2 mmol) in N,N-dimethylformamide (100 ml) was heated at 70 °C for 24 hours. The reaction mixture was partitioned between ether and water, and the organic layer was washed with water and brine successively, dried over anhydrous sodium sulfate,

and concentrated under reduced pressure. The residue was purified by silica gel column chromatography (eluent: 10-20% ethyl acetate-hexane) to afford (R)-N-benzyl-2-azido-2-(4-chlorophenyl)ethylamine (12.7 g, 83.8%) as a yellow oil.  $^{1}$ H NMR (CDCl<sub>3</sub>)  $\delta$ : 2.81(1H, dd, J=12.5, 5.1Hz), 2.89(1H, dd, J=12.5, 8.5Hz), 3.82(2H, s),4.64(1H, dd, J=8.5, 5.1Hz),7.23-7.36(9H, m).

A solution of (R)-N-benzyl-2-azido-2-(4-chlorophenyl)ethylamine (12.7 g, 44.1 mmol) in tetrahydrofuran (176 mL) was treated with triphenylphosphine (13.9 g, 52.9 mmol) at room temperature. After addition of water (20 ml), the reaction mixture was heated at 60 °C for 1 h. The reaction mixture was condensed, and partitioned between ether and 1 N hydrochloric acid. The aqueous layer was treated with 1 N aqueous sodium hydroxide solution until the solution became basic. The resulting solution was extracted with dichlromethane thoroughly. The combined organic layer was washed with water, dried over anhydrous sodium sulfate, filtered, and concentrated under reduced pressure. The residue was used for the next reaction without further purification.

The residue was heated with diethyl oxalate (18 ml, 132 mmol) at 120 °C for 1.5 h. The resulting white precipitate was washed with ether and collected to afford (R)-1-benzyl-5-(4-chlorophenyl)-2,3-dioxopiperazine (11.4 g, 82.2%). <sup>1</sup>H NMR (CDCl<sub>3</sub>)  $\delta$ : 3.46(1H, dd, J=12.9, 8.1Hz), 3.60(1H, dd, J=12.9, 3.8Hz), 4.48(1H, d, J=14.7Hz), 4.79(1H, d, J=14.7Hz), 4.80(1H, dd, J=8.9, 3.8Hz), 6.83(1H, s), 7.13(4H, m), 7.27(5H, m).

To a suspension of (R)-1-benzyl-5-(4-chlorophenyl)-2,3-dioxopiperazine (11.4 g, 36.3 mmol) in tetrahydrofuran (300 ml) was added borane-tetrahydrofuran complex (181 mL, 1.0 M solution in tetrahydrofuran, 181 mmol) at room temperature. After stirring for 24 hours, the reaction mixture was quenched with methanol (50 ml) at 0 °C, and concentrated under reduced pressure. The residue was treated with 10% aqueous sodium hydroxide solution (300 ml) and heated at 100 °C for 2 hours. After cooling at room temperature, the mixture was extracted with dichloromethane thoroughly. The combined organic layer was dried over anhydrous sodium sulfated, filtered, and concentrated under reduced pressure. The residue was used for the next reaction without further purification.

To a solution of the residue and triethylamine (7.58 ml, 54.4 mmol) in

dichloromethane (150 ml) was added di-tert-butyl dicarbonate (9.49 g, 43.5 mmol) at room temperature. After stirring for 45 min, the resulting mixture was partitioned between dichloromethane and water, dried over anhydrous sodium sulfate, filtered, and concentrated under reduced pressure. The residue was purified by silica gel column chromatography (eluent: 10-20% ethyl acetate-hexane) to afford (R)-1-benzyl-4-(tert-butoxycarbonyl)-3-(4-chlorophenyl)piperazine (11.6 g, 82.8%) as an oil.

<sup>1</sup>H NMR (CDCl<sub>3</sub>)  $\delta$ : 1.43(9H, s), 2.16(1H, dt, J=4.4, 11.7Hz), 2.40(1H, dd, J=4.4, 11.7Hz), 2.78(1H, dd, J=4.4, 11.7Hz), 2.98(1H, dt, J=4.4, 11.7Hz), 3.20(1H, d, J=12.8Hz), 3.42(1H, d, J=12.9Hz), 3.57(1H, d, J=12.9Hz), 3.89(1H, d, J=12.8Hz), 5.17(1H, s), 7.24-7.36(9H, m).

To a solution of (R)-1-benzyl-4-(tert-butoxycarbonyl)-3-(4-chlorophenyl)piperazine (11.6 g, 30.1 mmol) in 1,2-dichloroethane (80 ml) was added 1-chloroethyl chloroformate (4.91 ml, 45.1 mmol) at room temperature. Upon disappearance of the starting material, the reaction mixture was concentrated under reduced pressure. The residue was then dissolved in methanol (100 ml) and refluxed for 30 min. The resulting white precipitate was filtered and washed with methanol to afford (R)-2-(4-chlorophenyl)piperazine dihydrochloride, which was liberated with aqueous sodium hydroxide solution, and extracted with dichloromethane to afford (R)-2-(4-chlorophenyl)piperazine (3.04 g, 51.4%) as white solid.

<sup>1</sup>H NMR (CDCl<sub>8</sub>)  $\delta$ :2.65(1H, dd, J=12.0, 10.5Hz), 2.82-3.04(4H, m), 3.09(1H, d, J=12.6Hz), 3.73(1H, dd, J=10.1, 2.7Hz), 7.29(4H, m)

The filtrate was concentrated under reduced pressure and partitioned between ether and 1 N hydrochloric acid. The aqueous layer was neutralized with 1 N aqueous sodium hydroxide solution, and extracted with dichloromethane thoroughly. The combined organic extracts were dried over anhydrous sodium sulfate, filtered, and concentrated under reduced pressure. The residue was purified after Boc-protection (Boc<sub>2</sub>O, Et<sub>3</sub>N, CH<sub>2</sub>Cl<sub>2</sub>) to furnish (R)-1,4-di(tert-butoxycarbonyl)-2-(4-chlorophenyl)piperazine (2.70 g, 22.6%) as pale yellow solid.

<sup>1</sup>H NMR (CDCl<sub>8</sub>)  $\delta$ : 1.43(9H, s), 1.46(9H, s), 2.96(2H, m), 3.32(1H, dd, J=13.8,

4.2Hz), 3.74(1H, m), 3.94(1H, d, J=11.4Hz), 4.40(1H, d, J=13.2Hz),5.23(1H, s),7.25(2H, m)

To a suspension of (R)-2-(4-chlorophenyl)piperazine dihydrochloride (1.09 g, 4.05 mmol) in tetrahydrofuran (24 ml) was added triethylamine (2.82 ml, 20.3 mmol). After stirring for 15 min at room temperature, 2-chloro-3-methyl-6-(4pyridyl)-3H-pyrimidin-4-one (748 mg, 3.38 mmol) was added portionwise. Upon disappearance of the chloropyrimidone, the reaction mixture was condensed under reduced pressure. The residue was partitioned between saturated aqueous sodium bicarbonate solution and dichloromethane. The organic layer was dried over anhydrous sodium sulfate, filtered, and concentrated under reduced pressure to give pale yellow solid, which was recrystallized from ethanol to afford (R)-2-(2-(4-chlorophenyl)piperazin-4-yl)-3-methyl-6-(4-pyridyl)-pyrimidin-4-one (998 mg, 77.4%) as white crystals. The enantiomer excess was determined by HPLC (>99% ee). The crystals were converted into its dihydrochloride salt. <sup>1</sup>H NMR (DMSO-d<sub>6</sub>)  $\delta: 3.40(3H, m), 3.46(3H, s), 3.62(1H, dd, J=12.0, 13.2Hz),$ 3.72(1H, m), 3.92(1H, t, J=15.5Hz), 4.68(1H, t, J=10.1Hz), 7.18(1H, s), 7.58(2H, d, J=8.6Hz), 7.83(2H, d, J=8.6Hz), 8.57(2H, d, J=6.6Hz), 9.01(2H, d, J=6.6Hz), 10.20(1H, d, J=7.8Hz), 10.76(1H, br s)MS: 382(M+H)  $[\alpha]_{D^{24}} = +62.2$  ° (c 1.00, H<sub>2</sub>O)

Example 6: Synthesis of (S)-2-(2-(4-chlorophenyl)piperazin-4-yl)-3-methyl-6-(4-pyridyl)-pyrimidin-4-one

(S)-isomer was prepared same as above by using (R)-2-methyl-CBS-oxazaborolidine instead of (S)-2-methyl-CBS-oxazaborolidine. 

1H NMR (DMSO-d<sub>6</sub>)  $\delta$ : 3.40 (3H, m), 3.45 (3H, s), 3.53-3.96 (3H, m), 4.68 (1H, t, J = 13.5Hz), 7.10 (1H, s), 7.60 (2H, d, J=8.3Hz), 7.76 (2H, d, J=8.3Hz), 8.38 (1H, br s), 8.91 (1H, d, J=4.8Hz), 9.88 (1H, br s), 10.31 (1H, br s)

MS: 382(M+H)

[ $\alpha$ ] $_{D^{24}}$  = -63.3 ° (c 1.00, H<sub>2</sub>O)

The compounds in the following table were prepared in the same manner as

the methods described above. The compound numbers in the following table correspond to those shown in the above-described table of preferred compounds.

Table 3

NO	NMR	Exact-MS
A19	2.51-2.89(4H, m), 3.31-3.34(4H, m), 3.39(3H,s), 3.56(2H, s), 6.80(1H, s), 7.25-7.31(1H, m), 7.31-7.36(4H, m), 7.98(2H, dd, J=1.5, 4.8 Hz), 8.68(2H, dd, J=1.5, 4.5 Hz)(DMSO-d6)	362
A25	3.32-3.34(4H, m), 3.46(3H, s), 3.48-3.51(4H, m), 6.80-6.85(1H, m), 6.84(1H, s), 7.01(2H, d, J=8.0 Hz), 7.23-7.28(2H, m), 8.00(2H, dd, J=1.3, 4.6 Hz), 8.70(2H, dd, J=1.5, 4.5 Hz)(DMSO-d6)	348
A156	3.47(3H,s), 3.51-3.60(4H, m), 3.62-3.71(4H, m), 6.85(1H, s), 7.41-7.49(1H, m), 7.56-7.61(1H, m), 8.02(2H, dd, J=1.5, 4.5 Hz), 8.09(1H, d, J=8.1 Hz), 8.16(1H, d, J=8.1 Hz), 8.70(2H, dd, J=1.5, 4.8 Hz)(DMSO-d6)	405
A289	1.11-1.28(3H, m), 2.98-3.16(1H, m), 3.28-3.41(1H, m), 3.39(3H, s), 3.54-3.80(3H, m), 3.88-3.99(1H, m), 4.08-4.26(4H, m), 4.32-4.45(1H, m), 7.13(1H, s), 7.37-7.53(5H, m), 8.45(2H, d, J=5.8 Hz), 8.96(2H, d, J=6.0 Hz) (DMSO-d6)	434
A361	3.44(3H,s), 3.54-3.95(6H,m), 4.64(1H,brs), 7.11(1H,s), 7.42-7.51(3H,m), 7.74(2H,d,J=6.6Hz), 8.46(2H,d,J=5.7Hz), 8.94(2H,d,J=5.7Hz), 9.98(1H,brs), 10.46(1H, brs) (DMSO-d6).	348
A364	(DMSO-d6): 3.41-3.76(4H, m), 3.48(3H, s), 3.89-4.01(2H, m), 4.96(1H, m), 7.16(1H, s), 7.33-7.58(3H, m), 8.11(1H, dd, J=7.2, 7.2Hz), 8.52(2H, d, J=6.6Hz), 8.97(2H, d, J=6.6Hz), 10.04(1H, m), 10.66(1H, m).	366
A365	3.43(s, 3H), 3.51-3.96(m, 6H), 4.70(m, 1H), 7.00(s, 1H), 7.25(m, 1H), 7.54(m, 2H), 7.60(m, 1H), 8.20(d, J=5.7Hz, 2H), 8.80 (d, J=5.7Hz, 2H)(CDCl3)	366
· A366	2.27-2.85(1H, m), 2.94-3.08(3H, m), 3.43(3H,s), 3.59-3.67(2H, m), 3.94-3.97(1H, m), 6.81(1H, s), 7.19(2H, t, J=8.9 Hz), 7.50-7.55(2H, m), 7.96(2H, dd, J=1.6, 4.5 Hz), 8.68(2H, dd, J=1.5, 4.6 Hz)(DMSO-d6)	366
A366 (HCI)	3.35-3.50(2H, m), 3.46(3H, s), 3.58-3.75(2H, m), 3.86-3.97(2H, m), 4.68(1H, t, J=9.3 Hz), 7.15(1H, s), 7.35(2H, t, J=9.0 Hz), 7.82-7.87(2H, m), 8.48(2H, d, J=6.6 Hz), 8.96(2H, d, J=6.3 Hz), 9.55-10.08(1H, m), 10.54-10.70(1H, m)(DMSO-d6)	366

	(CDCl3):2.81(1H,dd,J=10.4,12.5Hz),	
A369	3.18-3.40(3H,m), 3.50-3.80(5H,m),	
	4.50(1H,dd,J=2.5,10.1Hz), 6.67(1H,s),	382
7.000	7.20-7.45(3H,m), 7.74(1H,dd,J=1.9,7.6Hz),	
	7.81(2H,dd,J=1.4,4.6Hz),	
	8.70(2H,dd,J=1.4,4.6Hz).	
	(CDCl3):3.01(1H,dd,J=10.4,12.5Hz),	
	3.10-3.30(3H,m), 3.50-3.80(5H,m),	
A370	4.04(1H,dd,J=2.7,10.8Hz), 6.67(1H,s),	382
	7.20-7.45(4H,m), 7.50(1H,s),	
	7.80(2H,dd,J=1.5,4.8Hz),	
	8.71(2H,dd, J=1.5,5.1Hz). 3.44(3H,s), 3.44-3.71(7H,m), 3.90(2H,m),	
	4.66(1H,brs), 7.11(1H,s),	
0.074	7.55(2H,d,J=8.4Hz), 7.78(2H,d,J=8.4Hz),	
A371	8.50(2H,d,J=5.7Hz), 8.95(2H,d,J=5.7Hz),	382
	10.13(1H,brs), 10.60(1H,brs)(DMSO-d6)	
	10.13(1H,D15), 10.00(1H,D15)(D1013O-00)	
1		
4070	(DMSO-d6):3.45(3H,s), 3.50-4.20(6H,m),	400
A376	4.66(1H,br s), 7.12(1H,s), 7.72(4H,s),	426
	8.44(2H,d,J=6.6Hz), 8.94(2H,d,J=6.6Hz),	
	10.00(1H,br s), 10.05(1H,br s).	
	3.37-3.93(6H, m), 3.48(3H, s), 3.87(3H, s),	
1	4.89-4.95(1H, m), 7.04-7.12(2H, m),	
A391	7.17(1H, d, J=8.5 Hz), 7.45-7.51(1H, m),	378
	7.75-7.81(1H, m), 8.29-8.38(2H, m),	
1	8.83-8.91(2H, m), 9.66-9.77(1H, m),	
<del> </del>	9.91-10.10(1H, m)(DMSO) (DMSO-d6):3.30-3.58(5H,m),	
	3.58-3.80(2H,m), 3.81(3H,s),	
	3.85-4.00(2H,m), 4.58-4.75(1H,m),	
A392	7.03(1H,dd,J=1.8,8.1Hz), 7.11(1H, s),	378
A392	7.26(1H,d,J=7.8Hz), 7.35-7.50(2H,m),	370
	8.41(2H,d,J=5.7Hz), 8.92(2H,d,J=6.0Hz),	
İ	9.80-10.00(1H,brd), 10.30-10.60(1H,brd).	
	3.40-3.43(5H,m), 3.51-3.63(2H,m),	
	3.78(3H,s), 3.93(2H,m),4.58(1H,br),	
A393	7.02-7.06(3H,m), 7.64(2H,d,J=8.7Hz),	378
7.000	8.34(2H,d,J=6.3Hz), 8.88(2H,d,J=8.7Hz),	0,0
1	9.76(1H,br), 10.16(1H,br)(DMSO-d6)	
····	1.30(3H, t, J=6.9 Hz), 3.38-3.54(1H, m),	
	3.49(3H, s), 3.65-3.79(1H, m), 3.84-3.98(2H,	
ļ	m), 4.02-4.18(2H, m), 4.84(1H, t, J=10.5 Hz),	
A396	7.04-7.16(2H, m), 7.15(1H, s), 7.39-7.45(1H,	392
AUSU	m), 7.89(1H, d, J=6.6 Hz), 8.49(2H, d, J=6.3	
	Hz), 8.95(2H, d, J=6.6 Hz), 9.92(1H, d, J=9.3	
	Hz), 10.51-10.64(1H, m)(DMSO-d6)	
	(DMSO-d6):3.64(2H,m),	
	3.94(2H,t,J=11.4Hz), 4.02-4.40(5H,m),	1
	4.78(1H,t,J=10.4Hz), 7.06(1H,s),	
A406	7.98(2H,d,J=8.3Hz), 8.01(2H,d,J=8.3Hz),	373
	8.23(1H,dd,J=1.2,5.1Hz),	
	9.02(1H,d,J=5.1Hz), 9.31(1H,d,J=1.2Hz),	1
l	10.03(1H,d,J=8.7Hz), 10.57(1H,s).	<u> </u>

A433	(CDCI3):2.00(4H,m), 3.03(1H,dd,J=10.8,12.0Hz), 3.21(3H,m), 3.29(4H,m), 3.57(3H,s), 3.62(2H,m), 3.90(1H,dd,J=2.7,10.8Hz), 6.57(2H,d,J=8.7Hz), 6.66(1H,s), 7.29(2H,d,J=8.7Hz), 7.80(2H,d,J=4.8Hz), 8.70(2H,d,J=4.8Hz).	417
A439	(CDCl3):3.02(1H,dd,J=10.7,12.4Hz), 3.18(7H,m), 3.55(3H,s), 3.62(2H,m), 3.87(4H,m), 3.96(1H,dd,J=2.5,11.1Hz), 6.66(1H,S), 6.93(2H,d,J=8.7Hz), 7.36(2H,d,J=8.7Hz), 7.79(2H,d,J=4.5Hz), 8.70(2H,d,J=4.5Hz).	434
A442	(CDCl3):2.36(3H,s), 2.59(4H,m), 3.02(1H,t,J=11.6Hz), 3.22(7H,m), 3.55(3H,s), 3.63(2H,m), 3.94(1H,d,J=10.5Hz), 6.66(1H,s), 6.93(2H,d,J=8.7Hz), 7.34(2H,d,J=8.7Hz), 7.80(2H,d,J=4.5Hz), 8.70(2H,d,J=4.5Hz).	.446
A463	3.41-3.54(3H, m), 3.48(3H, s), 3.69-3.73(1H, m), 3.78(3H, s), 3.82(3H, s), 3.86-3.93(2H, m), 4.89(1H, t, J=10.5 Hz), 6.97-7.01(1H, m), 7.08(1H, d, J=9.0 Hz), 7.15(1H, s), 7.66(1H, d, J=3.0 Hz), 8.51(2H, d, J=6.3 Hz), 8.96(2H, d, J=6.3 Hz), 9.93(1H, d, J=9.0 Hz), 10.60-10.73(1H, m)(DMSO-d6)	408
A464	(DMSO-d6): 3.45(3H, s), 3.38-3.81(6H, m), 3.88(6H, s), 5.06(1H, m), 6.82(2H, d, J=8.7Hz), 7.04(1H, s), 7.44(1H, t, J=8.4Hz), 8.20(1H, m), 8.30(2H, d, J=6.3Hz), 8.87(2H, d, J=6.3Hz), 10.07(1H, m).	408
A468	3.40-3.50(4H, m), 3.47(3H, s), 3.83-3.94(2H, m), 3.88(3H, s), 4.81-4.91(1H, m), 6.92-6.99(1H, m), 7.07-7.10(1H, m), 7.12(1H, s), 7.79-7.91(1H, m), 8.30-8.40(2H, m), 8.85-8.92(2H, m), 9.70-9.79(1H, m), 10.02-10.23(1H, m)(DMSO)	396
A469/ A470	(DMSO-d6):3.38-3.60(6H,m), 3.60-3.80(1H,m):3.80-4.00(5H,m), 4.80-4.97(1H,m); 6.85-7.00(1H,m), 7.09(1H,dd,J=2.4,11.4Hz); 7.13(1H,s), 7.95(1H,dd,J=6.9,8.7Hz), 8.46(2H,d,J=6.6Hz); 8.94(2H,d,J=6.3Hz); 9.80-10.00(1H,brd); 10.35-10.60(1H,brd).	396
A472	3.36-4.00(6H, m), 3.46(3H, s), 3.94(3H, s), 4.94-5.02(1H, m), 6.96-7.01(1H, m), 7.05(1H, d, J=8.6 Hz), 7.14(1H, s), 7.49-7.58(1H, m), 8.44-8.50(2H, m), 8.52-8.64(1H, m), 8.96(2H, d, J=6.6 Hz), 10.49-10.60(1H, m)(DMSO)	396
A480	2.78(1H, dd, J=10.0, 12.1 Hz), 3.18-3.27(3H, m), 3.59(3H, s), 3.64-3.74(2H, m), 3.86(3H, s), 4.37(1H, dd, J=2.4, 10.1 Hz), 6.67(1H, s), 6.89(1H, d, J=2.1 Hz), 6.99(1H, dd, J=1.7, 8.0 Hz), 7.50(1H, d, J=8.2 Hz), 7.82(2H, dd, J=1.5, 4.8 Hz), 8.71(2H, dd, J=1.8, 4.5 Hz)(CDCI3)	412

A490 (2HCl)	3.35-3.94(6H, m), 3.49(3H, s), 4.71-4.80(1H, m), 7.02-7.11(1H, m), 7.18-7.28(2H, m), 7.98-8.10(1H, m), 8.31-8.48(2H, m), 8.87-8.97(2H, m), 9.79-9.92(1H, m), 10.18-10.39(1H, m) (DMSO)	380
A501	(CDCI3):2.77(1H,dd,J=10.2,12.0Hz), 3.15-3.35(3H,m), 3.50-3.80(5H,m), 3.84(3H,s), 4.39(1H,d,J=7.8Hz), 6.67(1H,s), 6.78(1H,d,J=8.8Hz), 7.39(1H,dd,J=2.4,8.7Hz), 7.71(1H,d,J=2.3Hz), 7.82(2H,d,J=6.0Hz), 8.71(2H,d,J=6.0Hz).	456
A510	(CDCl3): 1.98-2.05(4H, m), 2.85(1H, dd, J=12, 10.5Hz), 3.17-3.24(7H, m), 3.58(3H, s), 3.65-3.72(2H, m), 3.85(3H, s), 4.28(1H, dd, 10.5, 2.7Hz), 6.10(1H, d, J=2.1Hz), 6.18(1H, dd, J=8.7, 2.1Hz), 6.65(1H, s), 7.33(1H, d, J=8.4Hz), 7.83(2H, dd, J=4.5, 1.8Hz), 8.70(2H, dd, J=4.5, 1.5Hz).	447
A511	(CDCl3):1.90-2.05(4H,m), 2.93(1H,t,J=12.0Hz), 3.15-3.40(7H,m), 3.59(3H,s), 3.65-3.85(5H,m), 4.11(1H,dd,J=2.1,10.2Hz), 6.49(1H,dd,J=3.0,9.0Hz), 6.66(1H,s), 7.83(2H,dd,J=1.8,4.5Hz), 8.70(2H,dd,J=1.5,4.5Hz).	447
A516	(DMSO-d6):3.20-3.70(4H,m), 3.70(1H,m), 3.98(3H,s), 3.99(3H,s), 4.00(1H,m), 4.96(1H,d,J=10.2Hz), 7.01(1H,s), 7.03(2H,m), 8.26(2H,d,J=6.1Hz), 8.53(1H,s), 8.84(2H,d,J=6.1Hz), 10.25(1H,d,J=10.7Hz)	414
A525	(DMSO-d6):3.30-3.50(2H,m), 3.48(3H,s), 3.55-3.78(2H,m), 3.78(3H,s), 3.96(2H,d,J=13.5Hz), 4.69(1H,t,J=10.4Hz), 7.06(1H,t,J=7.4Hz), 7.12(1H,s), 7.14(1H,d,J=7.4Hz), 7.31(1H,d,J=7.4Hz), 7.39(1H,t,J=7.4Hz), 7.59(2H,d,J=8.3Hz), 7.77(2H,d,J=8.3Hz), 8.43(2H,d,J=6.5Hz), 8.93(2H,d,J=6.5Hz), 9.89(1H,d,J=8.7Hz), 10.34(1H,s).	454
A527	(DMSO-d6):3.40-4.10(9H,m), 3.81(3H,s), 4.69(1H,m), 7.05(1H,s), 7.05(2H,d,J=9.0Hz), 7.67(2H,d,J=9.0Hz), 7.75(4H,s), 8.27(2H,d,J=5.7Hz), 8.85(2H,d,J=5.7Hz), 9.75(1H,s), 10.04(1H,s).	454
A536	(DMSO-d6):3.40-3.60(2H,m), 3.47(3H,s), 3.68(2H,m), 3.95(2H,m), 4.71(1H,t,J=9.9Hz), 7.16(1H,s), 7.33(2H,t,J=8.85Hz), 7.78(6H,m), 8.50(2H,d,J=6.3Hz), 8.97(2H,d,J=6.3Hz), 10.02(1H,s), 10.50(1H,s).	443
A543	3.52(s, 3H), 3.57-4.10(m, 6H), 5.57(m, 1H), 7.02(s, 1H), 7.53-7.70(m, 2H), 8.06(d, J=7.2Hz, 2H), 8.21-8.34(m, 3H), 8.82(d, J=6.3Hz, 2H), 9.88-9.92(m, 1H), 10.58-10.61(m, 1H)(DMSO d6)	398

A544	3.41-3.59(2H, m), 3.49(3H, s), 3.68-3.76(2H, m), 3.97-4.02(2H, m), 4.78-4.89(1H, m), 7.15(1H, s), 7.58-7.63(2H, m), 7.89-8.07(4H, m), 8.30(1H, s), 8.49(2H, d, J=6.3 Hz), 8.95(2H, d, J=6.3 Hz), 10.17(1H, d, J=8.4 Hz), 10.57-10.70(1H, m)(DMSO-d6)	398
A619	(CDCl3): 2.98(1H, dd, J=12.6, 10.8Hz), 3.17-3.28(5H, m), 3.58(3H, s), 3.62(1H, m), 3.79(1H, m), 4.26(1H, dd, 10.5, 2.7Hz), 4.62(2H, m), 6.66(1H, s), 6.88(1H, t, J=7.5Hz), 7.16(1H, d, J=7.2Hz), 7.27(1H, m), 7.84(2H, d, J=6.0), 8.70(2H, dd, J=4.8, 1.2Hz).	390
A626	3.33-3.41(4H, m), 3.42(3H, s), 3.47-3.87(4H, m), 6.84(1H, s), 7.44-7.49(5H, m), 7.99(2H, dd, J=1.5, 4.5 Hz), 8.69(2H, dd, J=1.4, 4.8 Hz)(DMSO-d6)	376
A649	3.44(3H, s),3.37-4.04(9H, m),4.67(1H, d,J=9.6Hz),7.10(1H, s),7.45-7.55(3H, m),7.83(2H, d,J=6.0Hz),8.47(2H, d,J=6.6Hz),8.95(2H, d,J=6.6Hz),12.15(1H, brs)(DMSO-d6)	362
A756	(CDCl3):2.50-2.61(1H,m), 2.80-2.95(1H,m), 3.05-3.20(1H,m), 3.25-3.40(1H,m), 3.50-3.60(1H,m), 3.57(3H,s), 3.65-3.75(1H,m), 3.75-3.80(1H,m), 3.85(3H,s), 6.60-6.80(3H,m), 7.47(1H,dd,J=7.2,8.4Hz), 7.82(2H,dd,J=1.5,4.5Hz), 8.71(2H,dd,J=1.5,4.5Hz).	410
A757/ A758	(DMSO-d6):2.54(3H,s), 3.40-3.79(3H,m), 3.46(3H,s), 3.80-4.10(6H,m), 4.83-5.10(1H,m), 6.90-7.05(1H,m), 7.08(1H,s), 7.13(1H,dd,J=2.7,11.4Hz), 8.00-8.25(1H,brd), 8.37(2H,d,J=6.3Hz), 8.91(2H,d,J=6.6Hz), 11.80-12.20(1H,brd).	410
A831	2.55(s, 3H), 3.51(s, 3H), 3.67-3.82(m, 4H), 4.04-4.08(m, 2H), 5.64(m, 1H), 7.05(s, 1H), 7.59-7.72(m, 3H), 8.06-8.11(m, 2H), 8.35(d, J=6.6Hz, 2H), 8.41(d, J=7.8Hz, 1H), 8.49 (d, J=6.9Hz, 1H), 8.84(d, J=6.6Hz, 2H)(DMSO d6)	412
A1016	(DMSO-d6):3.15-3.35(1H,m), 3.38-3.60(4H,m), 3.75-4.15(8H,m), 4.18-4.25(1H,m), 4.90-5.20(1H,m), 7.00-7.20(3H,m), 7.30-7.55(6H,m), 8.50-8.70(3H,m), 9.00(2H,d,J=6.3Hz).	. 486
A1276	(CDCl3):1.80-2.42(3H, m), 3.08-3.39(4H, m), 3.40-3.62(1H, m), 3.65-4.23(6.8H, m), 4.63-4.90(0.6H, m), 5.40-5.62(0.7H, m), 5.80-6.00(0.1H, m), 6.52-6.78(3H, m), 6.90-7.2(1H, m), 7.68-7.90(2H, m), 8.64-8.80(2H, m)	438

	1.48(3H, s), 1.57(3H, s), 3.50(3H, s),	
	3.51-3.66(2H, m), 3.72-3.76(1H, m),	
	3.90(3H, s), 3.99(1H, d, J=13.4 Hz),	
A1649	5.15-5.23(1H, m), 7.08-7.12(2H, m),	400
A1049	7.18(1H, d, J=8.6 Hz), 7.46-7.49(1H, m),	406
[	8.04-8.11(1H, m), 8.37-8.45(2H, m),	
	8.89-8.97(2H, m), 9.49-9.60(1H, m),	
	9.95-10.11(1H, m)(DMSO)	
	1.16-1,28(1H, m), 1.50-1.64(1H, m),	
1	1.70-1.82(2H, m), 1.90-2.01(1H, m),	
	2.58(2H, d, J=7.3 Hz), 2.64-2.72(1H, m),	
B13	2.89-2.97(1H, m), 3.28(3H, s), 3.57-3.67(2H,	361
-	m), 6.93(1H, s), 7.20-7.35(5H, m), 8.26(2H,	301
1	d, J=5.7 Hz), 8.87(2H, d, J=5.9	
	Hz)(DMSO-d6)	
	1.75-2.16(4H, m), 2.96-3.08(3H, m, 3.55(3H,	<del></del>
	11.73-2.10(4H, HI), 2.30-3.00(3H, M, 3.55(3H,	
B16	s), 3.69-3.79(2H, m), 6.66(1H, s),	347
1	7.26-7.40(5H, m), 7.81(2H, d, $J = 6.0 \text{ Hz}$ ),	
	8.70(2H, d, J = 6.0 Hz) (CDCI3)	
	1.76-1.99(5H, m), 2.97-3.10(2H, m),	
	3.75(1H, d, J=12.4 Hz), 6.81(1H, s),	
B17	7.18-7.24(2H, m), 7.28-7.35(1H, m),	365
	7.47(1H, t, J-7.1 Hz), 7.98(2H, d, J-5.8 Hz),	
	8.68(2H, d, J=5.8 Hz)(DMSO-d6)	
	1.86-2.14(4H, m), 2.94-3.03(3H, m),	
B19	3.55(3H, s), 3.68-3.75(2H, m), 6.66(1H, s),	205
513	7.05(2H, m), $7.23(2H, m)$ , $7.80(2H, d, J = 6.3)$	365
	Hz), $8.70(2H, d, J = 6.3 Hz)(CDCI3)$	
i	1.75-2.08(4H, m), 2.80(1H, m), 3.03(1H, m),	
	3.42(3H, s), 3.77(2H, m), 3.85(3H, s),	
B33	6.65(1H, s), 6.89-7.00(2H, m), 7.22-7.28(2H,	<b>37</b> 7
1	m), $7.82(2H, d, J = 6.0 Hz)$ , $8.70(2H, d, J = 1)$	
	6.0 Hz) (CDCl3)	
	1.73-1.83(4H, m), 2.90-3.02(3H, m),	
}	3.42(3H, s), 3.67-3.81(2H, m), 3.74(3H, s),	
B35	6.80(1H, s), 6.91(2H, d, J=8.7 Hz), 7.27(2H,	<b>37</b> 7
i	d, J=8.5 Hz), 7.97(2H, d, J=5.9 Hz), 8.69(2H,	
	d, J=5.7 Hz)(DMSO-d6)	
	1.69-1.90(7H, m), 1.94-2.00(1H, m),	
	2.59-2.68(4H, m), 2.92-3.02(3H, m),	
B43	3.43(3H, s), 3.69-3.80(4H, m), 6.59(3H, s),	430
	6.79(1H, s), 7.29-7.36(4H, m), 7.96(2H, d,	
	J=5.9 Hz), 8.68(2H, d, J=5.1 Hz)(DMSO-d6)	
	(CDCl3): 1.95-2.09(3H, m), 2.39(1H, m),	
	3.15(1H, m), 3.45(1H, dd, J=12.9, 10.8Hz),	
	3.57(3H, s), 3.61-3.72(2H, m), 4.08(1H, m),	
B46	6.67(1H, s), 7.32(1H, m), 7.58-7.60(2H, m),	388
1	7.74(1H, d, J=7.8Hz), 7.80(2H, dd, J=4.5,	
	1.5Hz), 8.69(2H, dd, J=4.5, 1.5Hz).	
<del></del>	(GDCl3): 1.90-2.06(3H, m), 2.36(1H, m),	
j	3.14(1H, m), 3.42(1H, m), 3.57(3H, s),	
B47	3.61-3.71(2H, m), 4.06(1H, m), 6.68(1H, s),	
	7.09(1H, m), 7.28(1H, m), 7.68(1H, dd,	406
	J=8.8, 5.1Hz), 7.79(2H, d, J=4.7Hz),	
	8.69(2H, d, J=5.9Hz).	
L	[ 0.03(20, u, 3-3.302).	<u> </u>

B48	1.90-2.10(3H, m), 2.32-2.44(1H, m), 3.11-3.20(1H, m), 3.45(1H, dd, J=10.5, 12.6 Hz), 3.57(3H, s), 3.61-3.72(2H, m), 4.08(1H, d, J=11.1 Hz), 6.67(1H, s), 7.30-7.35(1H, m), 7.56-7.62(2H, m), 7.74(1H, d, J=13.8 Hz), 7.80(2H, dd, J=1.8, 4.5 Hz), 8.70(2H, dd, J=1.8, 4.8 Hz)(CDCI3)	388
B49	1.91-2.09(3H, m), 2.37-2.42(1H, m), 3.12-3.19(1H, m), 3.45(1H, dd, J=10.8, 12.9 Hz), 3.57(3H, s), 3.60-3.72(2H, m), 4.08(1H, d, J=11.1 Hz), 6.67(1H, s), 7.30-7.35(1H, m), 7.54-7.62(2H, m), 7.75(1H, d, J=8.1 Hz), 7.80(2H, dd, J=1.5, 4.5 Hz), 8.70(2H, dd, J=1.8, 4.5 Hz)(CDCI3)	388
B50	1.59-1.67(1H, m), 1.72-1.81(1H, m), 2.08(1H, dt, J=3.4, 12.7 Hz), 2.23-2.40(1H, m), 3.06-3.14(1H, m), 3.41-3.54(2H, m), 3.42(3H, s), 3.93(1H, d, J=14.0 Hz), 7.02(1H, s), 7.24-7.29(1H, m), 7.34-7.39(2H, m), 7.56-7.59(2H, m), 8.55(2H, d, J=6.6 Hz), 8.98(2H, d, J=6.5 Hz)(DMSO-d6)	363
B80	2.21-2.36(4H, m), 3.19-3.31(2H, m), 3.46(3H, s), 3.88(2H, d, J=13.2 Hz), 6.86(1H, s), 7.38-7.42(1H, m), 7.46-7.51(2H, m), 7.58-7.64(2H, m), 8.01(2H, d, J=5.1 Hz), 8.70(2H, d, J=5.1 Hz)(DMSO-d6)	372
B122	1.44(2H, m), 1.75-1.83(3H, m), 2.63(2H, d, J = 6.9 Hz), 2.90(2H, m), 3.51(3H, s), 3.64(2H, m), 6.65(1H, s), 7.17-7.34(5H, m), 7.80(2H, d, J = 6.3 Hz) (CDCI3)	361
B123	1.44-2.16(5H, m), 2.86-2.97(2H, m), 3.49(3H, s), 3.62(1H, m), 3.72(1H, m), 4.48(1H, d, J = 7.2 Hz), 6.64(1H, s), 7.07(2H, m), 7.32(2H, m), 7.79(2H, d, J = 6.3 Hz), 8.69(2H, d, J = 6.3 Hz) (CDCI3)	395
B124	1.38-1.60(3H, m), 1.78(1H, m), 2.16(1H, m), 2.79-2.94(2H, m), 3.20(3H, s), 3.49(3H, s), 3.59(1H, m), 3.69(1H, m), 3.88(1H, d, J = 7.5 Hz, 1H), 6.64(1H, s), 7.08(2H, m), 7.25(2H, m), 7.79(2H, d, J = 6.0 Hz), 8.70(2H, d, J = 6.0 Hz) (CDCl3)	409
B127	1.87-2.06(4H, m), 2.79(1H, m), 3.10(2H, m), 3.57(3H, s), 3.78(2H, m), 6.68(1H, s), 7.23-7.29(3H, m), 7.34(2H, m), 7.84(2H, d, J = 6.0 Hz), 8.72(2H, d, J = 6.0 Hz) (CDCI3)	347
B130	1.81-2.03(4H, m), 2.78(1H, m), 3.09(2H, m), 3.57(3H, s), 3.79(2H, m), 6.69(1H, s), 7.03(2H, m), 7.23(2H, m), 7.84(2H, d, J = 5.4 Hz), 8.72(2H, br s) (CDCI3)	365
B134	1.78-1.95(4H, m), 2.80-2.91(1H, m), 2.96-3.09(2H, m), 3.45(3H, s), 3.81(2H, d, J=13.1 Hz), 6.80(1H, s), 7.33(1H, dd, J=2.0, 8.3 Hz), 7.56-7.60(2H, m), 7.99(2H, dd, J=1.6, 4.5 Hz), 8.69(2H, dd, J=1.5, 4.5 Hz)(DMSO-d6)	415

	1.82-2.02(4H, m), 3.09-3.27(3H, m),	
	3.57(3H, s), 3.79(2H, m), 3.86(3H, s),	
B145	6.67(1H, s), 6.89-6.99(2H, m), 7.21-7.26(2H,	377
	m), 7.84(2H, d, $J = 6.0$ Hz), 8.72(2H, d, $J = 6.0$ Hz)	
	6.0 Hz) (CDCl3) 1.85-2.07(2H,m), 2.17-2.30(2H,m),	
	2.91-3.10(1H,m), 3.10-3.24(2H,m),	
	3.57(3H,s), 3.71-3.88(2H,m), 6.69(1H,s),	
B157	6.99-7.06(1H,m), 7.21(1H,dd,J=2.1,8.7Hz),	405
	7.45(1H,s), 7.49-7.65(1H,m),	403
	7.83(2H,dd,J=1.8,4.5Hz),	
	8.72(2H,dd,J=1.2,4.8Hz)(CDCl3)	
	2.22-2.32(4H, m), 3.22(2H, m), 3.37(1H, m),	
	3.58(3H, s), 3.82(2H, m, 6.71(1H, s),	
B158	7.10(1H, m), 7.29(1H, m), 7.67(1H, m),	406
	7.83(2H, d, $J = 6.3$ Hz), 8.72(2H, d, $J = 6.3$	
,	Hz) (CDCl3)	
	2.19-2.26(4H, m), 3.21(2H, m), 3.35(1H, m),	
DAEO	3.59(3H, s), 3.82(2H, m), 6.70(1H, s),	405
B159	6.95(1H, dt, $J = 9.0$ , 2.1 Hz), 7.13(1H, dd, $J = 9.0$ , 2.1 Hz), 7.74(4H, m), 7.85(2H, d, $J = 9.0$ )	405
	9.0, 2.1 Hz), 7.71(1H, m), 7.85(2H, d, J = 6.3 Hz), 8.72(2H, d, J = 6.3 Hz) (CDCI3)	
	2.13-2.34(2H,m), 2.34-2.43(2H,m),	
	3.10-3.38(3H,m), 3.57(3H,s),	•
D400	3.68-3.83(2H,m), 6.69(1H,s),	
B160	7.29-7.40(2H,m), 7.46-7.59(1H,m),	388
	7.64-7.78(1H,m), 7.80-7.78(2H,m),	
	8.72(2H,d,J=6.0Hz)(CDCl3)	
	2.19(2H, m), 2.38(2H, m), 3.18(2H, m),	
	3.39(1H, m), 3.58(3H, s), 3.80(2H, m),	
B161	6.70(1H, s), 7.39(1H, m), 7.50(1H, m),	404
	7.83(2H, d, J = 6.0 Hz), 7.89(1H, d, J = 7.2	
	Hz), 8.01(1H, d, J = 7.8 Hz), 8.73(2H, d, J =   6.0 Hz) (CDCl3)	
	1.96(2H, m), 2.88(2H, m), 3.15(2H, m),	
	3.60(3H, s), 3.85(2H, m), 4.63(1H, m),	
B162	6.73(1H, s), 7.13-7.23(3H, m), 7.46(1H, d, J	420
	= 7.5 Hz), 7.84(2H, d, J = 6.3 Hz), 8.73(2H, d,	
	J = 6.3 Hz)(CDCI3)	
	1.64(2H, m), 2.23(2H, m), 3.13(2H, m),	
D464	3.50(1H, m), 3.53(3H, s), 3.68(2H, m),	
B164	6.58(2H, m), 6.68(1H, s), 6.91(2H, m),	380
	7.81(2H, d, J = 6.0 Hz), 8.72(2H, d, J = 6.0 Hz) (CDCI3)	
	1.91-1.99(4H, m), 2.84(3H, s), 3.07(2H, m),	<del></del>
	3.55(3H, s), 3.77(2H, m), 3.84(1H, m),	_
B165	6.69(1H, s), 6.75-6.87(3H, m), 7.27(2H, m),	376
	7.82(2H, d, J = 6.3 Hz), 8.72(2H, d, J = 6.3	- · ·
	Hz) (CDCI3)	
	1.52(2H, m), 1.79(3H, s), 1.96(2H, m),	
B168	3.09(2H, m), 3.42(3H, s), 3.64(2H, m),	
	4.86(1H, m), 6.63(1H, s), 7.09-7.19(4H, m),	422
1	7.74(2H, d, J = 6.0 Hz), 8.70(2H, d, J = 6.0 Hz)	
ļ	Hz) (CDCl3)	
	1.86(1H, br s), 1.95(2H, m), 2.30(2H, m), 3.47-3.63(7H, m), 6.68(1H, s), 7.30-7.44(3H, s)	
B169	3.47-3.63(7H, m), 6.68(1H, s), 7.30-7.44(3H, m), 7.54(2H, d, J = 7.5 Hz), 7.84(2H, d, J =	363
	6.0 Hz), 8.71(2H, d, J = 6.0 Hz) (CDCl3)	
i		i

B201	2.20-2.31(4H, m), 3.20-3.29(2H, m), 3.46(3H, s), 3.87(2H, d, J=13.8 Hz), 6.86(1H, s), 7.29-7.35(2H, m), 7.64-7.69(2H, m), 8.01(2H, dd, J=1.5, 4.5 Hz), 8.70(2H, dd, J=1.5, 4.5 Hz)(DMSO-d6)	390
B227	2.16-2.25(2H, m), 2.48-2.58(2H, m), 3.14-3.21(2H, m), 3.40(3H, s), 3.41-3.50(2H, m), 6.79(1H, s), 7.28-7.33(1H, m), 7.39-7.46(4H, m), 7.97(2H, dd, J=1.5, 4.5 Hz), 8.68(2H, dd, J=1.5, 4.5 Hz)(DMSO-d6)	389

Test Example: Inhibitory activity of the medicament of the present invention against P-GS1 phosphorylation by bovine cerebral TPK1

A mixture containing 100 mM MES-sodium hydroxide (pH 6.5), 1 mM magnesium acetate, 0.5 mM EGTA, 5 mM  $\beta$ -mercaptoethanol, 0.02% Tween 20, 10% glycerol, 12  $\mu$  g/ml P-GS1, 41.7  $\mu$  M [ $\gamma$ -32P] ATP (68 kBq/ml), bovine cerebral TPK1 and a compound shown in Table (a final mixture contained 1.7% DMSO deriving from a solution of a test compound prepared in the presence of 10% DMSO) was used as a reaction system. The phosphorylation was started by adding ATP, and the reaction was conducted at 25°C for 2 hours, and then stopped by adding 21% perchloric acid on ice cooling. The reaction mixture was centrifuged at 12,000 rpm for 5 minutes and adsorbed on P81 paper (Whatmann), and then the paper was washed four times with 75 mM phosphoric acid, three times with water and once with acetone. The paper was dried, and the residual radioactivity was measured using a liquid scintillation counter. The results are shown in the table below. The test compound markedly inhibited the P-GS1 phosphorylation by TPK1. The results strongly suggest that the medicaments of the present invention inhibit the TPK1 activity, thereby suppress the A  $\beta$  neurotoxicity and the PHF formation, and that the medicaments of the present invention are effective for preventive and/or therapeutic treatment of Alzheimer disease and the above-mentioned diseases.

Table 4

Compound No.	IC50
A19	$0.319\mu\mathrm{M}$
A25	$0.065\mu\mathrm{M}$
A156	0.140 μ M
A289	2.01 μ M
A361	0.018 μ M
A364	11nM
A365	0.022 μ M
A366	18nM
A366 (HCl)	7.7nM
A369	9.6 nM
A370	5.4 nM
A371	0.0066 μ Μ
A376	22nM
A391	1.8nM
A392	11nM
A393	0.0047 μ Μ
A396	25nM
A406	15 nM
A433	21 nM
A439	13 nM
A442	2.0 nM
A463	2.2nM
A464	4.7nM
A468	0.9nM
A469, A470	54nM
A469, A470	0.40nM
A472	5.5nM
A480	0.31 nM
A490(2HC1)	24nM .
A501	1.5 nM
A510	2.5nM
A511	3.7 nM
A516	2.0nM
A525	32 nM
A527	5.0 nM
A536	1.6nM
A543	$0.0012\mu\mathrm{M}$
A544	4.6nM

A619     1.6 nM       A626     0.093 μ M       A649     0.855 μ M       A756     56nM       A757, A758     1181nM	
A649 0.855 μ M A756 56nM	
A756 56nM	
<u> </u>	
A757, A758 1181nM	
A757, A758 34nM	
A831 0.015 μ M	
A1016 1170 nM	
A1276 1674 nM	
A1649 7.7nM	
B13 0.013 μ M	
Β16 0.019 μ Μ	
B17 0.0053 μ M	
B19 0.024 μ M	
B33 0.0027 μ M	
B35 0.0075 μ M	
B43 0.0013 μ M	
B46 0.0018 μ M	
B47 0.0088 μ M	
B48 0.0098 μ M	
B49 0.00078 μ M	_
B50 0.351 μ M	
B80 0.23 μ M	_
B122 0.052 μ M	
B123 0.087 μ M	
B124 0.147 μ M	
B127 0.026 μ M	
B130 0.024 μ M	
B134 0.11 μ M	
B145 0.033 μ M	
B157 0.020 μ M	
B158 0.011 μ M	_
B159 . 0.0054 μ M	
B160 0.0099 μ M	
B161 0.0056 μ M	
B162 0.015 μ M	
B164 0.028 μ M	
B165 0.042 μ M	
B168 0.041 μ M	
B169 0.023 μ M	

B201	$0.27~\mu\mathrm{M}$
B227	$0.062~\mu$ M

# Formulation Example

#### (1) Tablets

The ingredients below were mixed by an ordinary method and compressed by using a conventional apparatus.

Compound of Example 1 30 mg
Crystalline cellulose 60 mg
Corn starch 100 mg
Lactose 200 mg
Magnesium stearate 4 mg

#### (2) Soft capsules

The ingredients below were mixed by an ordinary method and filled in soft capsules.

Compound of Example 1 30 mg

Olive oil 300 mg

Lecithin 20 mg

### Effect of Invention

The compounds of the present invention have TPK1 inhibitory activity and are useful as an active ingredient of a medicament for preventive and/or therapeutic treatment of diseases caused by abnormal advance of TPK1 such as neurodegenerative diseases (e.g. Alzheimer disease) and the above-mentioned diseases.

# 【書類名】外国語要約書

# ABSTRACT

A pyrimidone derivative having tau protein kinase 1 inhibitory activity which is represented by formula (I) or a salt thereof, or a solvate thereof or a hydrate thereof:

$$(X)_{m} \xrightarrow{N}_{R} O$$

$$(Y)_{n} \xrightarrow{R}$$

$$(X)_{m} \xrightarrow{N}_{R} O$$

wherein R represents a C1-C12 alkyl group; the ring of:

represents piperazine ring or piperidine ring; each X independently represents a C1-Cs alkyl group, an optionally partially hydrogenated C6-C10 aryl ring, an indan ring or the like; m represents an integer of 1 to 3; each Y independently represents a halogen atom, a hydroxy group, a cyano group, a C1-C6 alkyl group or the like; n represents an integer of 0 to 8; when X and Y or two Y groups are attached on the same carbon atom, they may combine to each other to form a C2-C6 alkylene group.

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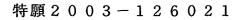
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